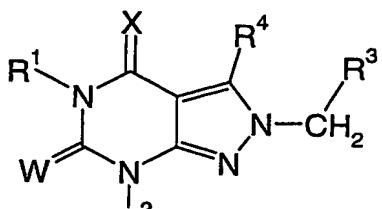


Claims:

1. A compound having the structural formula (I):



5

(I)

wherein,

X is S, O, or NR²⁰, provided that when W is O, then X is not O,

X and the double bond to which it is attached can be replaced with 2 hydrogen atoms,

10 W is S, O, or NR²⁰; provided that when X is O, then W is not O;

R¹ is H, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, hydroxy, amino, or optionally substituted heterocycle;

15 R² is H, optionally substituted alkyl, optionally substituted alkylcycloalkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, or optionally substituted heterocycle;

20 R³ is a monocyclic or bicyclic, saturated or unsaturated, ring system comprising 0, 1, 2 or 3 heteroatoms independently selected from N, O, or S, the ring being substituted by 0, 1, 2 or 3 substituents selected from =O, halogen, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, nitro, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -NR^aR^a, -NHC(=O)R^a, N=NR^a, aminocarbonyl, phenyl, benzyl; or R³ is represented by -Het, -Het-Het, R⁵, -R⁵-Het, -Het-R⁵, -Het-O-R⁵, -R⁵-R⁵, -R⁵-OR⁵;

25 R⁴ is a monocyclic or bicyclic, saturated or unsaturated, ring system, or a vicinal-fused derivative thereof, which may contain from 5 to 12, preferably 5 to 10, ring atoms, 0, 1, 2, 3

or 4 of which are heteroatoms independently selected from N, O, or S, the ring system being substituted by 0, 1, 2 or 3 substituents selected from B(OH)₂, vicinal -OCH₂CH₂O-, vicinal -OC₁₋₂haloalkylO-, vicinal -OCH₂O-, vicinal -CH₂OCH₂O-, =O, halogen, -R^bOR^a, -SR^a, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -

5 -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -NR^aR^a, -NHC(=O)R^a, -NHC(=O)OR^a, N=NR^a, NO₂, -C(=O)NR^aR^a, -C(=O)NR^aOR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)_nR^a), -

10 C(=O)NR^a(R^bHet), -C(=O)OR^a, -OC(=O)R^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -C(=O)R^bNR^aR^a, -C(=O)NR^aR^a, -C(=NCN)R^a, -S(=O)₂NR^aR^a, -NR^aS(=O)₂R^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), -S(=O)₂NR^a(R^bC(=O)OR^a), aminocarbonyl, phenyl, benzyl; or R⁴ is represented by -(CH₂)_nR⁵-

15 Het, -(CH₂)_nR^d, -Het, -Het-Het, R⁵, -R⁵-Het, -Het-R⁵, -Het-OR⁵, R⁵-R⁵, or -R⁵-OR⁵; or R⁴ is represented by C₁₋₆alkyl, -NC₁₋₆alkyl, or -N(C₁₋₆alkyl)₂ wherein the C₁₋₆alkyl, -NC₁₋₆alkyl, -N(C₁₋₆alkyl) are substituted by 0, 1 or 2 substituents selected from R^a, OR^a, halogen or phenyl wherein R⁴ is not -(CH₂)₂CH₃, -(CH₂)₂CH₂OH, -(CH₂)₂CO₂H, or -(CH₂)₂CO₂C₁₋₆alkyl wherein z is 1,2,3,4,5, or 6;

R⁵ is independently at each instance, phenyl substituted by 0, 1, 2, or 3 groups selected from halogen, C₁₋₆haloalkyl, -OC₁₋₆haloalkyl, C₁₋₆alkyl, -CN, nitro, -OR^a, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a,

20 -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -R^bOR^a, -SR^a, -C(=O)NR^aR^a, -C(=O)NR^aOR^a, -C(=O)NR^aR^bNR^aR^a, -C(=O)NR^aR^bOR^a, -C(=O)NR^aR^bS(=O)_nR^a, -C(=O)NR^aR^bHet, -C(=O)OR^a, -OC(=O)R^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -C(=O)R^bNR^aR^a, -C(=O)NR^aR^a, -C(=NCN)R^a, -S(=O)₂NR^aR^a, -NR^aS(=O)₂R^a, -S(=O)₂NR^aR^bC(=O)NR^aR^a, or -S(=O)₂NR^aR^bC(=O)OR^a;

R²⁰ is, independently at each instance, H, -CN, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a;

R^a is, independently at each instance, H, C₁₋₆alkyl, -C(=O)C₁₋₄alkyl, C₁₋₄haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

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R^b is, independently at each instance, C_{1-6} alkyl, $-C(=O)C_{1-4}$ alkyl, C_{1-4} haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

R^c is C_{1-6} alkyl, C_{1-4} haloalkyl, phenyl or benzyl;

5 R^d is phenyl substituted by 0, 1 or 2 groups selected from -CN, halogen, nitro, C_{1-6} alkyl, C_{1-4} haloalkyl, -OH, -OR^a, -NR^aR^a, -S(=O)_nR^c, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, -OC(=O)R^a, B(OH)₂, vicinyl -OCH₂CH₂O-, vicinyl -OC₁₋₂haloalkylO-, vicinyl -OCH₂O-, vicinyl -CH₂OCH₂O-, phenyl, benzyl and a 5- or 6-membered ring, saturated or unsaturated heterocycle containing 1, 2, 3 or 4 heteroatoms independently selected from N, O, or S;

10 m is 1, 2 or 3;

 n is 0, 1 or 2;

When "optionally substituted" is used, it refers to at least one substituent selected from cyclopropyl, halogen, nitro, cyano, hydroxy, trifluoromethyl, amino, carboxy, carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} alkoxy, C_{1-4} alkanoyl, C_{1-4} alkanoyloxy, NH(C_{1-4} alkyl), N(C_{1-4} alkyl)₂, C_{1-4} alkanoylamino, (C_{1-4} alkanoyl)₂amino, N-(C_{1-4} alkyl)carbamoyl, N,N-(C_{1-4} alkyl)₂carbamoyl, (C_{1-4})S, (C_{1-4} alkyl)S(O), (C_{1-4} alkyl)S(O)₂, (C_{1-4}) alkoxycarbonyl, N-(C_{1-4} alkyl)sulfamoyl, N,N- C_{1-4} alkyl)sulfamoyl, C_{1-4} alkylsulfonylamino, and heterocyclic

20 or a pharmaceutically acceptable salt thereof.

2. A compound as recited in Claim 1 wherein:

R^1 is H, or C_{1-6} alkyl, or $-(CH_2)_n$ cycloalkyl or $-(CH_2)_{1-2}$ Het wherein C_{1-6} alkyl or $-(CH_2)_n$ cycloalkyl or $-(CH_2)_{1-2}$ Het is optionally substituted by 1, 2 or 3 substituents selected from Het, halogen, -CN, -OR^a, -NR^aR^a, -C(=O)OR^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl, -S(=O)_nR^c, -S(=O)_nNR^aR^a or -NR^aC(=O)C₁₋₄alkyl and n is 0, 1 or 2.

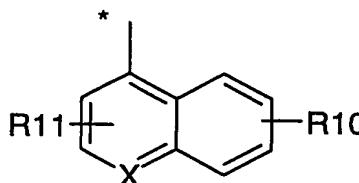
3. A compound as recited in Claim 1 wherein:

R^2 is $-(CH_2)_{1-3}$ cycloalkyl or $-C_{1-12}$ alkyl wherein $-(CH_2)_{1-3}$ cycloalkyl or $-C_{1-12}$ alkyl is optionally substituted with 0, 1, 2 or 3 substituents selected from Het, S(=O)_nR^c, -S(=O)_nNR^aR^a halogen, -CN, -OR^a, -NR^aR^a, -C(=O)OR^a, -C(=O)R^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl, or -NR^aC(=O)C₁₋₄alkyl and n is 0, 1 or 2.

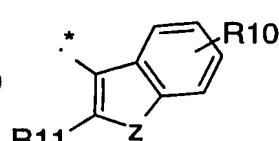
- 85 -

4. A compound as recited in Claim 1 wherein:

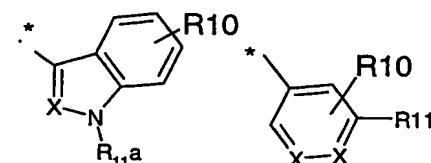
R^3 is selected from formulas (i), (ii), (iii) or (iv) set forth below:



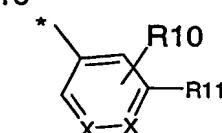
(i)



(ii)



(iii)



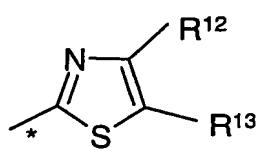
(iv)

wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R^{10} is at any position on the ring and R^{10} and R^{11} are independently at each instance H, R^a , halogen, -CN, nitro, OR^a , CF_3 , $-NR^aR^a$, $-C(=O)OR^a$, $C(=O)R^a$, $-C(=O)NR^aR^a$, $-OC(=O)C_{1-4}alkyl$, $-NR^aC(=O)C_{1-4}alkyl$ or $-S(=O)_nR^c$; and wherein $R^{11,a}$ is R^a , $-S(=O)_2NR^aR^a$ or $-S(=O)_nR^c$ and $n=1$ or 2.

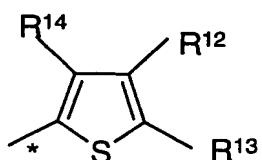
5. A compound as recited in Claim 1 wherein:

R^4 is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

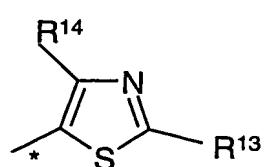
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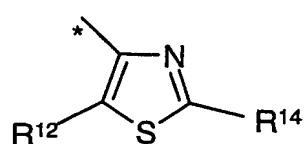
(a)



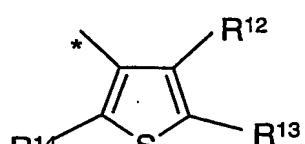
(b)



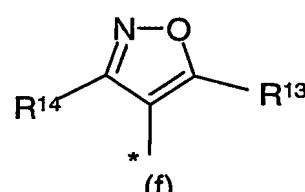
(c)



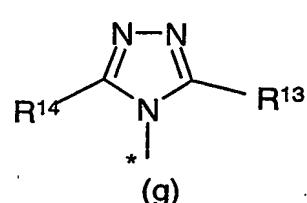
(d)



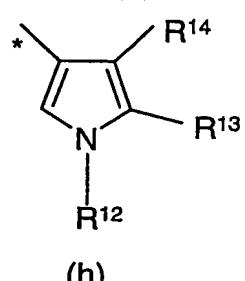
(e)



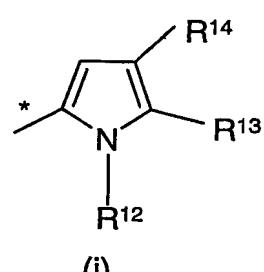
(f)



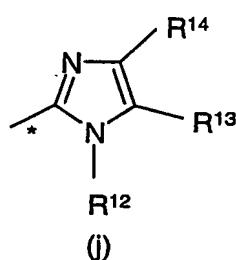
(g)



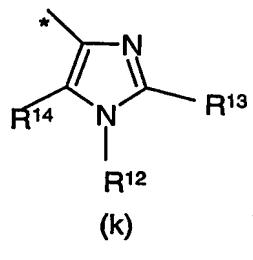
(h)



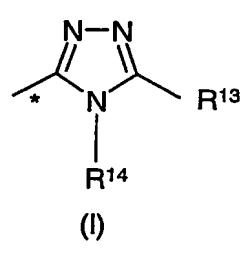
(i)



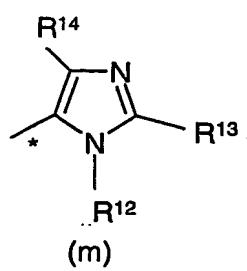
(j)



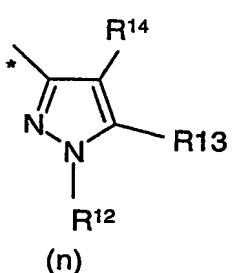
(k)



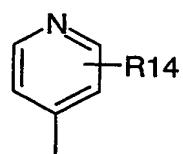
(l)



(m)

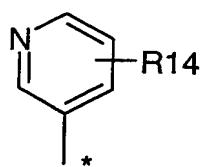


(n)

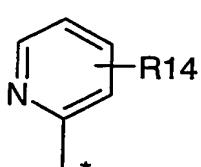


(o)

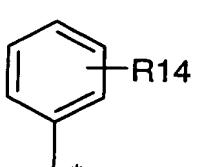
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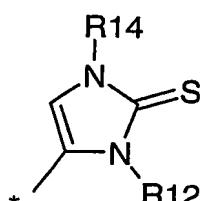
(p)



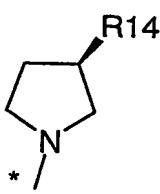
(q)



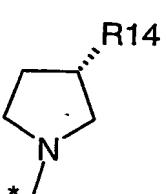
(r)



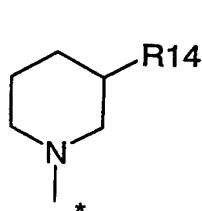
(s)



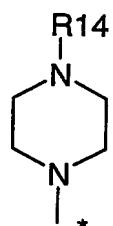
(t)



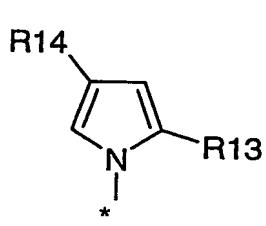
(u)



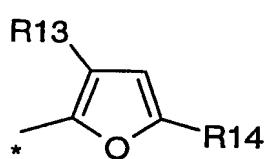
(v)



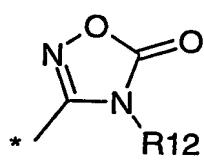
(w)



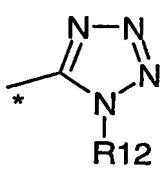
(x)



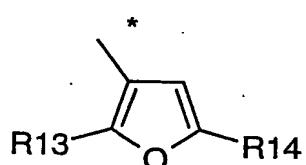
(y)



(z)



(aa)



wherein * is the location wherein R⁴ is attached to the ring system and wherein wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -

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C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a).

6. A compound as recited in Claim 1 wherein:

X is S, O, or NR²⁰, provided that when W is O, then X is not O; or X and the double bond to which it is attached can be 2 hydrogen atoms,

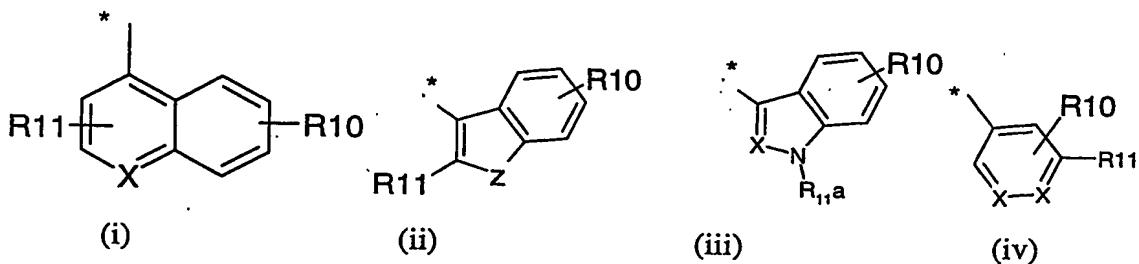
10 W is S, O, or NR²⁰; provided that when X is O, then W is not O;

R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a

15 R¹ is CH₃, CH₂CH₃, CH₂CN, CF₃, (CH₂)₂OH, cyclopropyl, isopropyl, CH₂CCH, (CH₂)₂N(CH₂)₂, (CH₂)₂N(C=NH)NH₂, -CH₂-2-pyridyl, -CH₂-3-pyridyl, -CH₂-4-pyridyl, -(CH₂)₂-1-imidazolyl, -(CH₂)₂-1-pyrazolyl, -(CH₂)₂-1-piperidyl, -(CH₂)_m-(1-methylpiperidin-4-yl), -CH₂-(1-methylpiperidin-3-yl), -(CH₂)₂-(morpholin-4-yl),

15 R² is -CH₂CH₂CH₃, -CH₂-cyclopropyl, -CH₂CH(CH₃)₂, -CH₂CH₂CH₂F, -CH₂-cyclobutyl, -CH₂C(CH₃)₃, -CH₂CH₂CH(CH₃)₂, -CH₂CF₃, -CH₂-methylphenyl, -CH₂-phenol, -CH₂-(3,5-dimethylisoxazol-4-yl), -CH₂-S-phenyl, -CH₂-phenylcarboxyl, or -CH₂SCF₃;

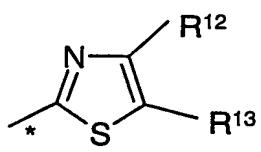
20 R³ is selected from formulas (i), (ii), (iii) or (iv) set forth below:



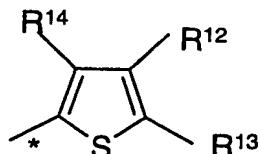
25 wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R¹⁰ is at any position on the ring and R¹⁰ and R¹¹ are independently at each instance H, R^a, halogen, -CN, nitro, OR^a, CF₃, -NR^aR^a, -C(=O)OR^a, -C(=O)R^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl, -NR^aC(=O)C₁₋₄alkyl or -S(=O)_nR^c; and wherein R^{11a} is R^a, -S(=O)₂NR^aR^a or -S(=O)₂NR^c and n=1 or 2.

30 R⁴ is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

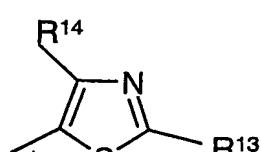
- 89 -



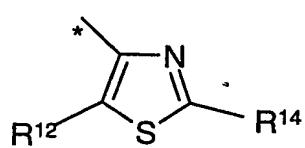
(a)



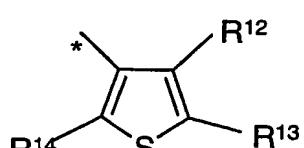
(b)



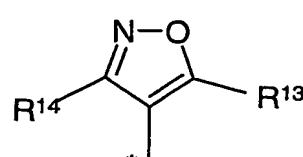
(c)



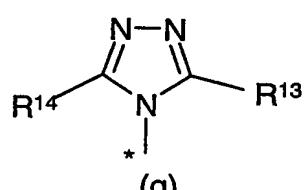
(d)



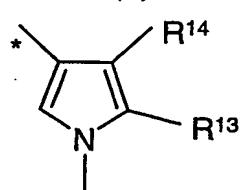
(e)



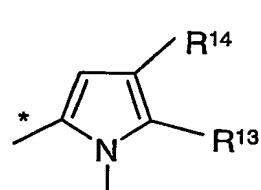
(f)



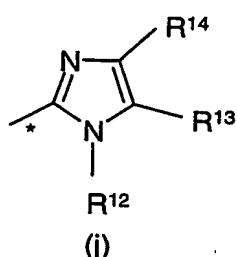
(g)



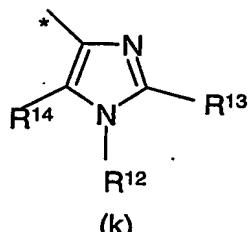
(h)



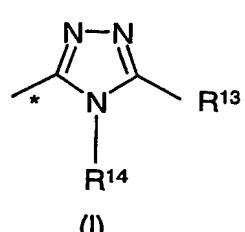
(i)



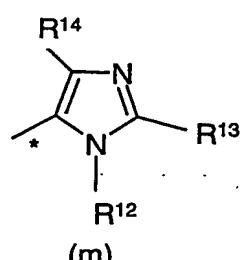
(j)



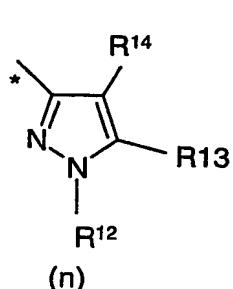
(k)



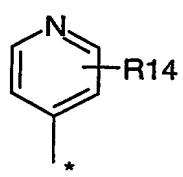
(l)



(m)

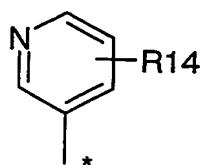


(n)

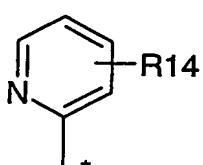


(o)

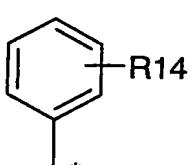
- 90 -



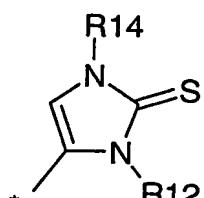
(p)



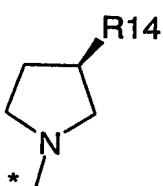
(q)



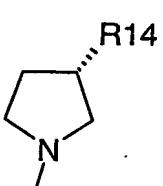
(r)



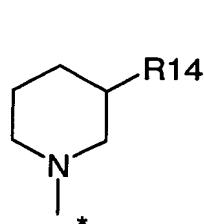
(s)



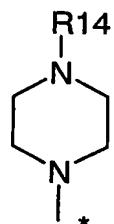
(t)



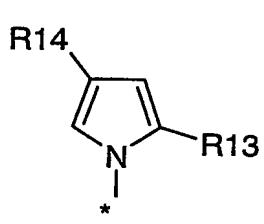
(u)



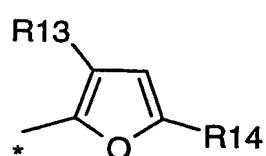
(v)



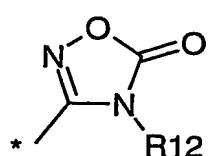
(w)



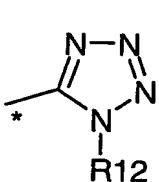
(x)



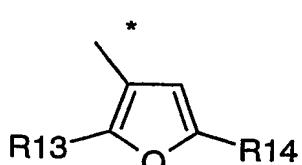
(y)



(z)



(aa)



(ab)

wherein * is the location wherein R⁴ is attached to the ring system and wherein
 wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -

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C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -

5 S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a.

7. A compound of formula (I) selected from:

5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-oxo-4-thioxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

10 5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-4-imino-5-methyl-6-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[(4Z)-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-4-(methylimino)-6-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

15 5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-imino-5-methyl-4-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[(2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-4-oxo-6-thioxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[(6Z)-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-6-(methylimino)-4-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

20 N-[(6Z)-2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-7-(cyclopropylmethyl)-5-methyl-4-oxo-2,4,5,7-tetrahydro-6H-pyrazolo[3,4-d]pyrimidin-6-ylidene]acetamide;

25 N-[(6Z)-2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-7-(cyclopropylmethyl)-5-methyl-4-oxo-2,4,5,7-tetrahydro-6H-pyrazolo[3,4-d]pyrimidin-6-ylidene]methanesulfonamide;

5-((6Z)-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-{[2-(dimethylamino)ethyl]imino}-5-methyl-4-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl)-1-methyl-1H-pyrrole-3-carbonitrile;

30 N-1~-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-7-(cyclopropylmethyl)-5-methyl-4-oxo-2,4,5,7-tetrahydro-6H-pyrazolo[3,4-d]pyrimidin-6-ylidene]-N~2~,N~2~-dimethylglycinamide;

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5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

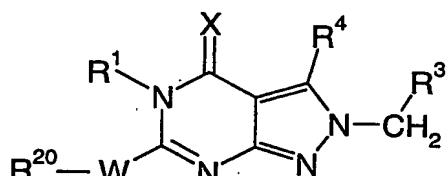
5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-6-oxo-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5 5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-3-(1-methyl-1H-imidazol-5-yl)-4-thioxo-2,4,5,7-tetrahydro-6H-pyrazolo[3,4-d]pyrimidin-6-one;

10 (4Z)-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-3-(1-methyl-1H-imidazol-5-yl)-4-(methylimino)-2,4,5,7-tetrahydro-6H-pyrazolo[3,4-d]pyrimidin-6-one.

8. A compound having the structural formula (II):



15

(II)

wherein,

X is S, O, or NR²⁰,

X and the double bond to which it is attached can be replaced with 2 hydrogen atoms,

20 W is S, O, or NR²¹;

R¹ is H, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, hydroxy, amino, or optionally substituted heterocycle, wherein the substitution is selected from cyclopropyl, halogen, nitro, cyano, hydroxy, trifluoromethyl, amino, carboxy,

25 carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C₁₋₄ alkyl, C₂₋₄ alkenyl, C₂₋₄ alkynyl, C₁₋₄ alkoxy, C₁₋₄ alkanoyl, C₁₋₄ alkanoyloxy, NH(C₁₋₄ alkyl), N(C₁₋₄ alkyl)₂, C₁₋₄ alkanoylamino, (C₁₋₄ alkanoyl)₂amino, N-(C₁₋₄ alkyl)carbamoyl, N,N-(C₁₋₄ alkyl)₂carbamoyl, (C₁₋₄)S, (C₁₋₄ alkyl)S(O), (C₁₋₄ alkyl)S(O)₂, (C₁₋₄) alkoxycarbonyl, N-(C₁₋₄ alkyl)sulfamoyl, 30 N,N-C₁₋₄ alkyl)sulfamoyl, C₁₋₄ alkylsulfonylamino, and heterocyclic;

R^3 is a monocyclic or bicyclic, saturated or unsaturated, ring system comprising 0, 1, 2 or 3 heteroatoms independently selected from N, O, or S, the ring being substituted by 0, 1, 2 or 3 substituents selected from =O, halogen, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, nitro,

5 -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a,
- $O(CH_2)_mNR^aR^a$, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a,
-S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -NR^aR^a, -NHC(=O)R^a, N=NR^a,
aminocarbonyl, phenyl, benzyl; or R³ is represented by -Het, -Het-Het, R⁵, -R⁵-Het, -Het-R⁵, -
Het-O-R⁵, -R⁵-R⁵, -R⁵-OR⁵;

10 R^4 is a monocyclic or bicyclic, saturated or unsaturated, ring system, or a vicinal-fused derivative thereof, which may contain from 5 to 12, preferably 5 to 10, ring atoms, 0, 1, 2, 3 or 4 of which are heteroatoms independently selected from N, O, or S, the ring system being substituted by 0, 1, 2 or 3 substituents selected from B(OH)₂, vicinal -OCH₂CH₂O-, vicinal -OC₁₋₂haloalkylO-, vicinal -OCH₂O-, vicinal -CH₂OCH₂O-, =O, halogen, -R^bOR^a, -SR^a, -OR^a,
C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -
O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -
S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -
S(CH₂)_mOR^a, -NR^aR^a, -NHC(=O)R^a, -NHC(=O)OR^a, N=NR^a, NO₂, -C(=O)NR^aR^a, -
C(=O)NR^aOR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)_nR^a), -
C(=O)NR^a(R^bHet), -C(=O)OR^a, -OC(=O)R^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -C(=O)R^bNR^aR^a,
-C(=NOR^a)R^a, -C(=NCN)R^a, -S(=O)₂NR^aR^a, -NR^aS(=O)₂R^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), -
S(=O)₂NR^a(R^bC(=O)OR^a), aminocarbonyl, phenyl, benzyl; or R⁴ is represented by -(CH₂)_nR⁵-
Het, -(CH₂)_nR^d, -Het, -Het-Het, R⁵, -R⁵-Het, -Het-R⁵, -Het-OR⁵, R⁵-R⁵, or -R⁵-OR⁵; or R⁴ is
represented by C₁₋₆alkyl, -NC₁₋₆alkyl, or -N(C₁₋₆alkyl)₂ wherein the C₁₋₆alkyl, -NC₁₋₆alkyl,
25 -N(C₁₋₆alkyl) are substituted by 0, 1 or 2 substituents selected from R^a, OR^a, halogen or
phenyl wherein R⁴ is not -(CH₂)₂CH₃, -(CH₂)₂CH₂OH, -(CH₂)₂CO₂H, or -(CH₂)₂CO₂C₁₋₆alkyl wherein z is 1,2,3,4,5, or 6;

30 R^5 is independently at each instance, phenyl substituted by 0, 1, 2, or 3 groups selected from halogen, C₁₋₆haloalkyl, -OC₁₋₆haloalkyl, C₁₋₆alkyl, -CN, nitro, -OR^a, -S(=O)_nR^c,
-O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a,
-O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a,
-S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -R^bOR^a, -SR^a, -C(=O)NR^aR^a, -
C(=O)NR^aOR^a, -C(=O)NR^aR^bNR^aR^a, -C(=O)NR^aR^bOR^a, -C(=O)NR^aR^bS(=O)_nR^a, -

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$C(=O)NR^aR^b$ Het, $-C(=O)OR^a$, $-OC(=O)R^a$, $-C(=O)OR^bNR^aR^a$, $-C(=O)R^a$, $-C(=O)R^bNR^aR^a$, $-C(=NOR^a)R^a$, $-C(=NCN)R^a$, $-S(=O)_2NR^aR^a$, $-NR^aS(=O)_2R^a$, $-S(=O)_2NR^aR^bC(=O)NR^aR^a$, or $-S(=O)_2NR^aR^bC(=O)OR^a$;

R^{20} is, independently at each instance, H, -CN, $-S(=O)_nR^c$, $-C(=O)R^a$, $-C(=O)NR^aR^a$,

5 $-C(=O)OR^a$, $-NR^aC(=O)R^a$, or $-OC(=O)R^a$, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle, wherein the substitution is selected from cyclopropyl, halogen, nitro, cyano, hydroxy, 10 trifluoromethyl, amino, carboxy, carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} alkoxy, C_{1-4} alkanoyl, C_{1-4} alkanoyloxy, $NH(C_{1-4})$ alkyl), $N(C_{1-4} \text{ alkyl})_2$, C_{1-4} alkanoylamino, $(C_{1-4} \text{ alkanoyl})_2$ amino, $N-(C_{1-4} \text{ alkyl})\text{carbamoyl}$, $N,N-(C_{1-4} \text{ alkyl})_2\text{carbamoyl}$, $(C_{1-4})S$, $(C_{1-4} \text{ alkyl})S(O)$, $(C_{1-4}\text{alkyl})S(O)_2$, $(C_{1-4})\text{alkoxycarbonyl}$, $N-(C_{1-4} \text{ alkyl})\text{sulfamoyl}$, $N,N-C_{1-4} \text{ alkyl})\text{sulfamoyl}$, C_{1-4} alkylsulfonylamino, and heterocyclic; 15

R^{21} is, independently at each instance, H, -CN, $-S(=O)_nR^c$, $-C(=O)R^a$, $-C(=O)NR^aR^a$,

$-C(=O)OR^a$, $-NR^aC(=O)R^a$, or $-OC(=O)R^a$; optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl,

20 optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle, wherein the substitution is selected from cyclopropyl, halogen, nitro, cyano, hydroxy, trifluoromethyl, amino, carboxy, carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} alkoxy, C_{1-4} alkanoyl, C_{1-4} alkanoyloxy, $NH(C_{1-4})$ alkyl), $N(C_{1-4} \text{ alkyl})_2$, C_{1-4} alkanoylamino, $(C_{1-4} \text{ alkanoyl})_2$ amino, $N-(C_{1-4} \text{ alkyl})\text{carbamoyl}$, 25 $N,N-(C_{1-4} \text{ alkyl})_2\text{carbamoyl}$, $(C_{1-4})S$, $(C_{1-4} \text{ alkyl})S(O)$, $(C_{1-4}\text{alkyl})S(O)_2$, $(C_{1-4})\text{alkoxycarbonyl}$, $N-(C_{1-4} \text{ alkyl})\text{sulfamoyl}$, $N,N-C_{1-4} \text{ alkyl})\text{sulfamoyl}$, C_{1-4} alkylsulfonylamino, and heterocyclic;

R^{20} and R^{21} and the N to which they are attached in combination can also form a 3 to

10 member N-linked saturated or unsaturated heterocycle having either 1, or 2 heteroatoms

30 independently selected from N, O, or S wherein the heterocycle is substituted with R^e ;

R^a is, independently at each instance, H, C_{1-6} alkyl, $-C(=O)C_{1-4}$ alkyl, C_{1-4} haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

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R^b is, independently at each instance, C₁₋₆alkyl, -C(=O)C₁₋₄alkyl, C₁₋₄haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

R^c is C_{1-6} alkyl, C_{1-4} haloalkyl, phenyl or benzyl;

5 R^d is phenyl substituted by 0, 1 or 2 groups selected from -CN, halogen, nitro,
C₁₋₆alkyl, C₁₋₄haloalkyl, -OH, -OR^c, -NR^aR^a, -S(=O)_nR^c, -C(=O)NR^aR^a, -C(=O)OR^a,
-NR^aC(=O)R^a, -OC(=O)R^a, B(OH)₂, vicinal -OCH₂CH₂O-, vicinal -OC₁₋₂haloalkylO-, vicinal
-OCH₂O-, vicinal -CH₂OCH₂O-, phenyl, benzyl and a 5- or 6-membered ring, saturated or
unsaturated heterocycle containing 1, 2, 3 or 4 heteroatoms independently selected from N, O,
10 or S:

10 or S;

R^e is independently at each instance, H, C_{1-6} alkyl, $-C(=O)C_{1-4}$ alkyl, C_{1-4} haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S:

m is 1, 2 or 3;

15 n is 0, 1 or 2;

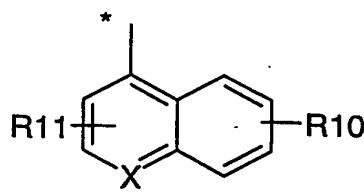
or a pharmaceutically acceptable salt thereof

9. A compound as recited in Claim 8 wherein:

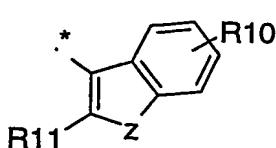
²⁰ R¹ is H, or C₁₋₆alkyl, or -(CH₂)_ncycloalkyl wherein C₁₋₆alkyl or -(CH₂)_ncycloalkyl is optionally substituted by 1, 2 or 3 substituents selected from Het, halogen, -CN, -OR^a, -NR^aR^a, -C(=O)OR^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl or -NR^aC(=O)C₁₋₄alkyl and n is 0, 1 or 2.

10. A compound as recited in Claim 8 wherein:

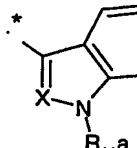
25 R^3 is selected from formulas (i), (ii), (iii) or (iv) set forth below:



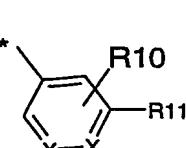
(i)



(ii)



(iii)



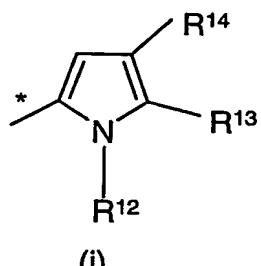
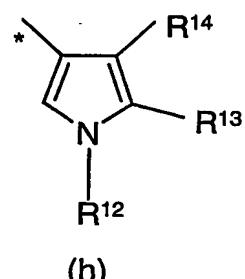
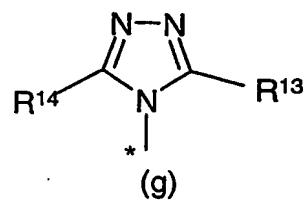
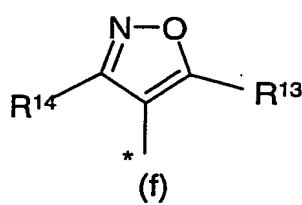
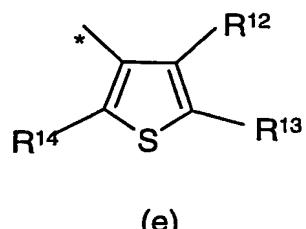
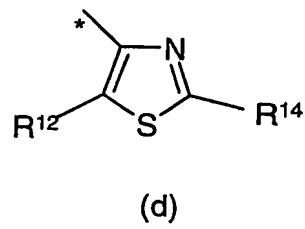
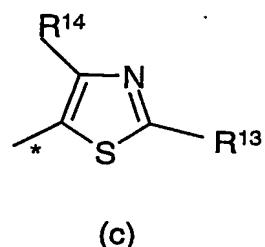
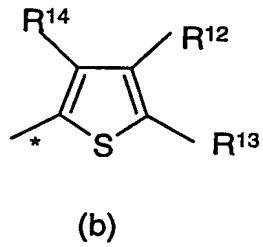
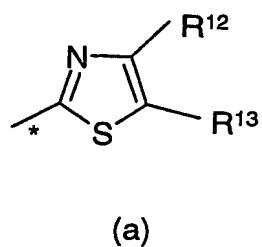
(iv)

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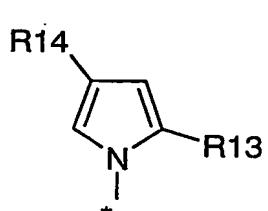
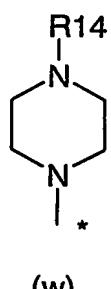
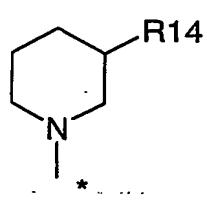
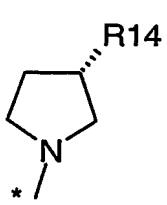
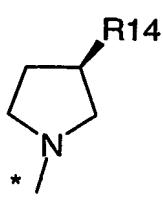
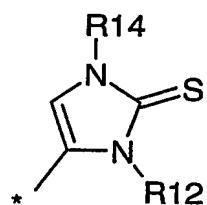
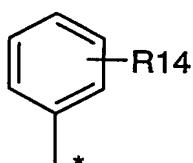
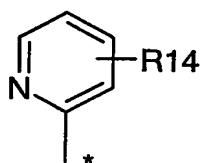
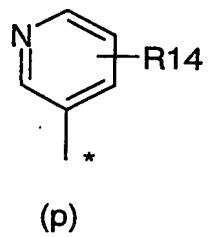
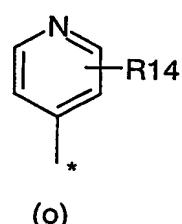
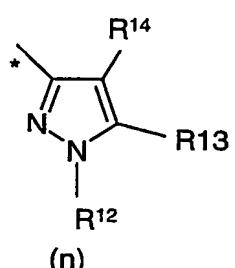
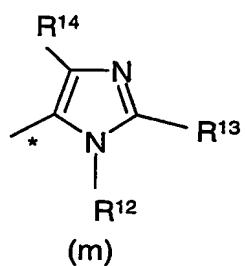
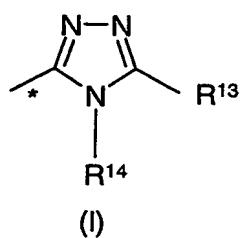
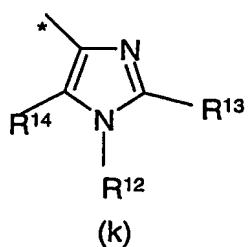
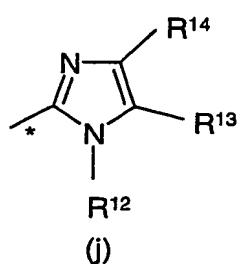
wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R¹⁰ is at any position on the ring and R¹⁰ and R¹¹ are independently at each instance H, R^a, halogen, -CN, nitro, OR^a, CF₃, -NR^aR^a, -C(=O)OR^a, -C(=O)R^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl, -NR^aC(=O)C₁₋₄alkyl or -S(=O)_nR^c; and wherein 5 R^{11a} is R^a, -S(=O)₂NR^aR^a or -S(=O)_nR^c and n=1 or 2.

11. A compound as recited in Claim 8 wherein:

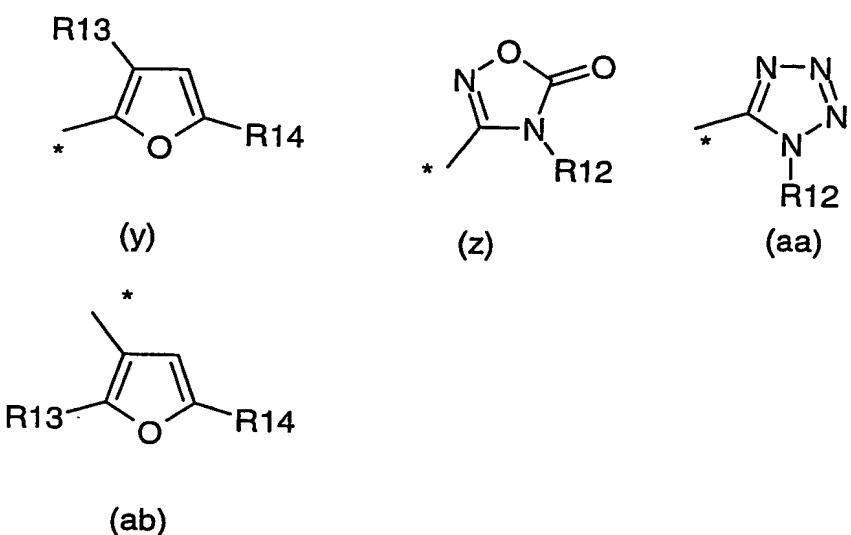
R⁴ is selected from formulas (a) to (z) or (aa) or (ab) set forth below:



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wherein * is the location wherein R⁴ is attached to the ring system and wherein
 wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁-alkyl, -CN,

5 NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a).

12. A compound as recited in Claim 8 wherein:

X is S, O, or NR²⁰; or X and the double bond to which it is attached can be 2 hydrogen atoms.

15 W is S, O, or NR²¹;

R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a,
-C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a;

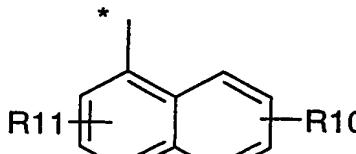
R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a,
-C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a;

20 R²⁰ and R²¹ and the N to which they are attached in combination can also form a 3 to
10 member N-linked saturated or unsaturated heterocycle having either 1, or 2 heteroatoms
independently selected from N, O, or S wherein the heterocycle is substituted with R^c.

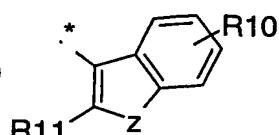
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R^1 is CH_3 , CH_2CH_3 , CH_2CN , CF_3 , $(CH_2)_2OH$, cyclopropyl, isopropyl, CH_2CCH , $(CH_2)_2N(CH_2)_2$, $(CH_2)_2N(C=NH)NH_2$, - CH_2 -2-pyridyl, - CH_2 -3-pyridyl, - CH_2 -4-pyridyl, - $(CH_2)_2$ -1-imidazolyl, - $(CH_2)_2$ -1-pyrazolyl, - $(CH_2)_2$ -1-piperidyl, - $(CH_2)_m$ -(1-methylpiperidin-4-yl), - CH_2 -(1-methylpiperidin-3-yl), - $(CH_2)_2$ -(morpholin-4-yl),

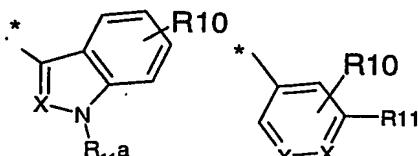
5 R^3 is selected from formulas (i), (ii), (iii) or (iv) set forth below:



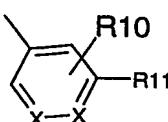
(i)



(ii)



(iii)

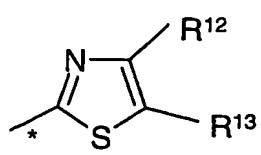


(iv)

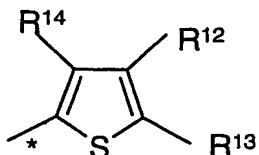
10 wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R^{10} is at any position on the ring and R^{10} and R^{11} are independently at each instance H, R^a , halogen, -CN, nitro, OR^a , CF_3 , $-NR^aR^a$, $C(=O)OR^a$, $-C(=O)R^a$, $-C(=O)NR^aR^a$, $-OC(=O)C_{1-4}alkyl$, $-NR^aC(=O)C_{1-4}alkyl$ or $-S(=O)_nR^c$; and wherein R^{11a} is R^a , $-S(=O)_2NR^aR^a$ or $-S(=O)_nR^c$ and $n=1$ or 2.

15 R^4 is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

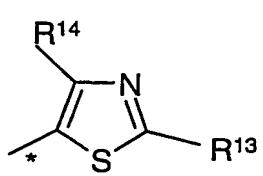
- 100 -



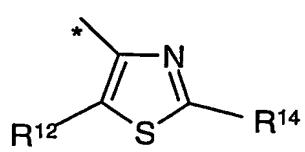
(a)



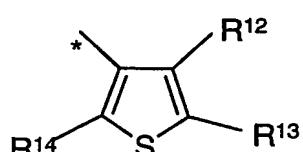
(b)



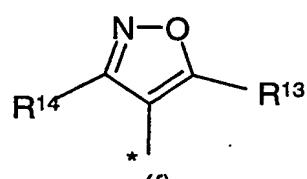
(c)



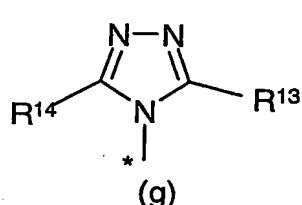
(d)



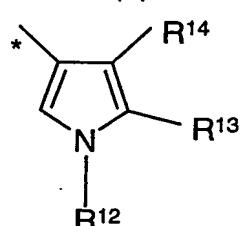
(e)



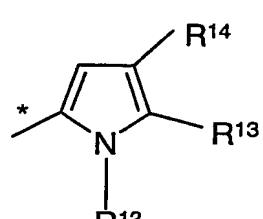
(f)



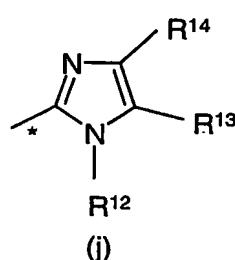
(g)



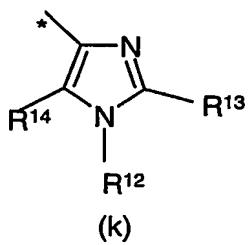
(h)



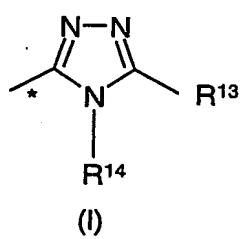
(i)



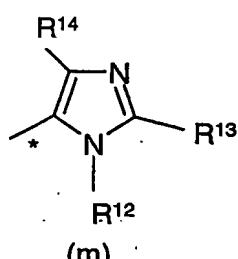
(j)



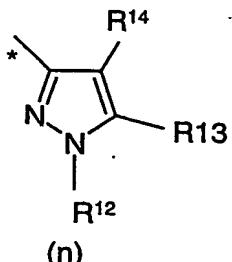
(k)



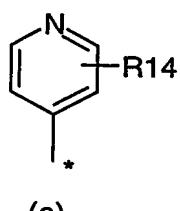
(l)



(m)

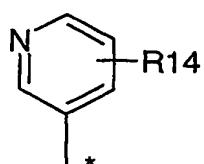


(n)

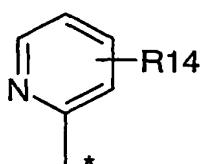


(o)

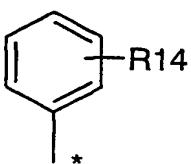
- 101 -



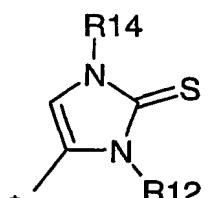
(p)



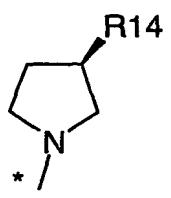
(q)



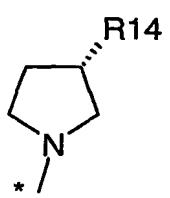
(r)



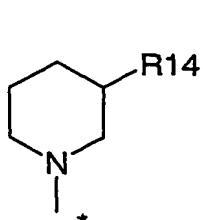
(s)



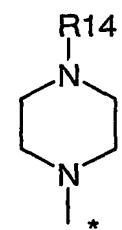
(t)



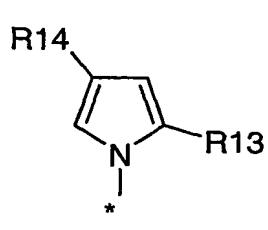
(u)



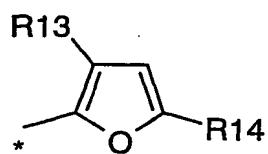
(v)



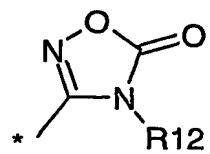
(w)



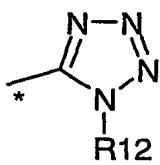
(x)



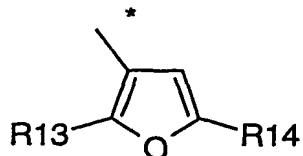
(y)



(z)



(aa)



5

(ab)

wherein * is the location wherein R⁴ is attached to the ring system and wherein
 wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -
 NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -

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C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a.

13. A compound of formula (II) selected from:

5-{6-amino-2-[(6-chloroquinolin-4-yl)methyl]-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile;

10 N-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-6-yl]-3-methylbutanamide
N,N-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-6-yl]-3-methylbutanamide;
N'-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-5-methyl-4-oxo-

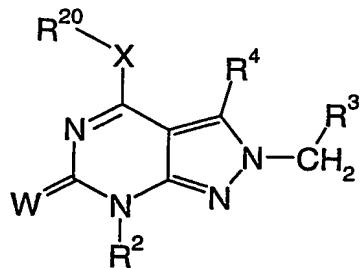
15 4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-6-yl]-N,N-dimethylimidoformamide;
5-{2-[(6-chloroquinolin-4-yl)methyl]-6-[(cyclopropylmethyl)(methyl)amino]-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile;
5-{2-[(6-chloroquinolin-4-yl)methyl]-6-[(cyclopropylmethyl)amino]-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile;

20 N-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-6-yl]propane-1-sulfonamide;
ethyl 2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-5-methyl-4-oxo-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-6-ylcarbamate;
N-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-5-methyl-4-oxo-

25 4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-6-yl]-N'-ethylurea;
5-[(4Z)-2-[(6-chloroquinolin-4-yl)methyl]-6-[(cyclopropylmethyl)amino]-5-methyl-4-(methylimino)-4,5-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;
5-[(4Z,6Z)-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-5-methyl-4,6-

30 bis(methylimino)-4,5,6,7-tetrahydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile.

14. A compound having the structural formula (III):



5 wherein,

X is S, O, NR²¹; or XR²⁰ is hydrogen;

W is S, O, or NR²⁰;

10 R² is H, optionally substituted alkyl, optionally substituted alkylcycloalkyl, optionally substituted alkylcycloalkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, or optionally substituted heterocycle;

15 R³ is a monocyclic or bicyclic, saturated or unsaturated, ring system comprising 0, 1, 2 or 3 heteroatoms independently selected from N, O, or S, the ring being substituted by 0, 1, 2 or 3 substituents selected from =O, halogen, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, nitro, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -NR^aR^a, -NHC(=O)R^a, N=NR^a, aminocarbonyl, phenyl, benzyl; or R³ is represented by -Het, -Het-Het, R⁵, -R⁵-Het, -Het-R⁵, -20 Het-O-R⁵, -R⁵-R⁵, -R⁵-OR⁵;

25 R⁴ is a monocyclic or bicyclic, saturated or unsaturated, ring system, or a vicinal-fused derivative thereof, which may contain from 5 to 12, preferably 5 to 10, ring atoms, 0, 1, 2, 3 or 4 of which are heteroatoms independently selected from N, O, or S, the ring system being substituted by 0, 1, 2 or 3 substituents selected from B(OH)₂, vicinal -OCH₂CH₂O-, vicinal -OC₁₋₂haloalkylO-, vicinal -OCH₂O-, vicinal -CH₂OCH₂O-, =O, halogen, -R^bOR^a, -SR^a, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mOR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a,

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$S(CH_2)_mOR^a$, $-NR^aR^a$, $-NHC(=O)R^a$, $-NHC(=O)OR^a$, $N=NR^a$, NO_2 , $-C(=O)NR^aR^a$, $-C(=O)NR^aOR^a$, $-C(=O)NR^a(R^bNR^aR^a)$, $-C(=O)NR^a(R^bOR^a)$, $-C(=O)NR^a(R^bS(=O)_nR^a)$, $-C(=O)NR^a(R^bHet)$, $-C(=O)OR^a$, $-OC(=O)R^a$, $-C(=O)OR^bNR^aR^a$, $-C(=O)R^a$, $-C(=O)R^bNR^aR^a$, $-C(=NOR^a)R^a$, $-C(=NCN)R^a$, $-S(=O)_2NR^aR^a$, $-NR^aS(=O)_2R^a$, $-S(=O)_2NR^a(R^bC(=O)NR^aR^a)$, $-S(=O)_2NR^a(R^bC(=O)OR^a)$, aminocarbonyl, phenyl, benzyl; or R^4 is represented by $-(CH_2)_nR^5$,

5 Het , $-(CH_2)_nR^d$, $-Het$, $-Het-Het$, R^5 , $-R^5-Het$, $-Het-R^5$, $-Het-OR^5$, R^5-R^5 , or $-R^5-OR^5$; or R^4 is represented by $C_{1-6}alky$, $-NC_{1-6}alkyl$, or $-N(C_{1-6}alkyl)_2$ wherein the $C_{1-6}alkyl$, $-NC_{1-6}alkyl$, $-N(C_{1-6}alkyl)$ are substituted by 0, 1 or 2 substituents selected from R^a , OR^a , halogen or phenyl wherein R^4 is not $-(CH_2)_zCH_3$, $-(CH_2)_zCH_2OH$, $-(CH_2)_zCO_2H$, or $-(CH_2)_zCO_2C_{1-}$

10 $alkyl$ wherein z is 1,2,3,4,5, or 6;

R^5 is independently at each instance, phenyl substituted by 0, 1, 2, or 3 groups selected from halogen, $C_{1-6}haloalkyl$, $-OC_{1-6}haloalkyl$, $C_{1-6}alkyl$, $-CN$, nitro, $-OR^a$, $-S(=O)_nR^c$, $-O(CH_2)_mHet$, $-O(CH_2)_mC(=O)Het$, $-O(CH_2)_mC(=O)NR^aR^a$, $-O(CH_2)_mC(=O)OR^a$, $-O(CH_2)_mNR^aR^a$, $-O(CH_2)_mOR^a$, $-S(CH_2)_mHet$, $-S(CH_2)_mC(=O)Het$, $-S(CH_2)_mC(=O)NR^aR^a$, $-S(CH_2)_mC(=O)OR^a$, $-S(CH_2)_mNR^aR^a$, $-S(CH_2)_mOR^a$, $-R^bOR^a$, $-SR^a$, $-C(=O)NR^aR^a$, $-C(=O)NR^aOR^a$, $-C(=O)NR^aR^bNR^aR^a$, $-C(=O)NR^aR^bOR^a$, $-C(=O)NR^aR^bS(=O)_nR^a$, $-C(=O)NR^aR^bHet$, $-C(=O)OR^a$, $-OC(=O)R^a$, $-C(=O)OR^bNR^aR^a$, $-C(=O)R^a$, $-C(=O)R^bNR^aR^a$, $-C(=NOR^a)R^a$, $-C(=NCN)R^a$, $-S(=O)_2NR^aR^a$, $-NR^aS(=O)_2R^a$, $-S(=O)_2NR^aR^bC(=O)NR^aR^a$, or $-S(=O)_2NR^aR^bC(=O)OR^a$;

20 R^{20} is, independently at each instance, H, $-CN$, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle, $-S(=O)_nR^c$, $-C(=O)R^a$, $-C(=O)NR^aR^a$, $-C(=O)OR^a$, $-NR^aC(=O)R^a$, or $-OC(=O)R^a$;

25 R^{21} is, independently at each instance, H, $-CN$, optionally substituted alkyl, optionally substituted alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle, $-S(=O)_nR^c$, $-C(=O)R^a$, $-C(=O)NR^aR^a$, $-C(=O)OR^a$, $-NR^aC(=O)R^a$, or $-OC(=O)R^a$; or

30 R^{20} and R^{21} and the N to which they are attached in combination can also form a 3 to 10 member N-linked saturated or unsaturated heterocycle having either 1, or 2 heteroatoms independently selected from N, O, or S wherein the heterocycle is substituted with R^e ;

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R^a is, independently at each instance, H, C_{1-6} alkyl, $-C(=O)C_{1-4}$ alkyl, C_{1-4} haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

R^b is, independently at each instance, C_{1-6} alkyl, $-C(=O)C_{1-4}$ alkyl, C_{1-4} haloalkyl,

5 phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

R^c is C_{1-6} alkyl, C_{1-4} haloalkyl, phenyl or benzyl;

R^d is phenyl substituted by 0, 1 or 2 groups selected from -CN, halogen, nitro,

10 C_{1-6} alkyl, C_{1-4} haloalkyl, -OH, -OR^c, -NR^aR^a, -S(=O)_nR^c, $-C(=O)NR^aR^a$, $-C(=O)OR^a$, -NR^aC(=O)R^a, -OC(=O)R^a, B(OH)₂, vicinyl -OCH₂CH₂O-, vicinyl -OC₁₋₂haloalkylO-, vicinyl -OCH₂O-, vicinyl -CH₂OCH₂O-, phenyl, benzyl and a 5- or 6-membered ring, saturated or unsaturated heterocycle containing 1, 2, 3 or 4 heteroatoms independently selected from N, O, or S;

R^e is independently at each instance, H, C_{1-6} alkyl, $-C(=O)C_{1-4}$ alkyl, C_{1-4} haloalkyl,

15 phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

m is 1, 2 or 3;

n is 0, 1 or 2;

When "optionally substituted" is used, it refers to at least one substituent selected from

20 cyclopropyl, halogen, nitro, cyano, hydroxy, trifluoromethyl, amino, carboxy, carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C_{1-4} alkyl, C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} alkoxy, C_{1-4} alkanoyl, C_{1-4} alkanoyloxy, NH(C_{1-4} alkyl), N(C_{1-4} alkyl)₂, C_{1-4} alkanoylamino, (C_{1-4} alkanoyl)₂amino, N-(C_{1-4} alkyl)carbamoyl, N,N-(C_{1-4} alkyl)₂carbamoyl, (C_{1-4})S, (C_{1-4} alkyl)S(O), (C_{1-4} alkyl)S(O)₂, (C_{1-4}) alkoxycarbonyl, N-(C_{1-4} alkyl)sulfamoyl, N,N- C_{1-4} alkyl)sulfamoyl, C_{1-4} alkylsulfonylamino, and heterocyclic

25 or a pharmaceutically acceptable salt thereof.

15. A compound as recited in Claim 14 wherein:

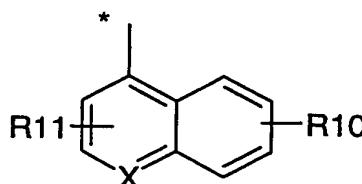
R^2 is $-(CH_2)_{1-3}$ cycloalkyl or $-C_{1-12}$ alkyl wherein $-(CH_2)_{1-3}$ cycloalkyl or $-C_{1-12}$ alkyl is

30 optionally substituted with 0, 1, 2 or 3 substituents selected from Het, S(=O)_nR^c, halogen, -CN, -OR^a, -NR^aR^a, -C(=O)OR^a, -C(=O)R^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl or -NR^aC(=O)C₁₋₄alkyl and n is 0, 1 or 2.

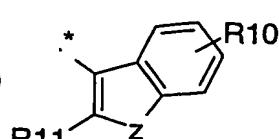
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16. A compound as recited in Claim 14 wherein:

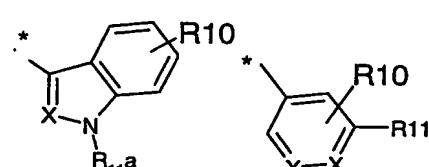
R^3 is selected from formulas (i), (ii), (iii) or (iv) set forth below:



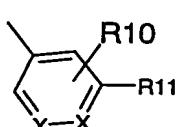
(i)



(ii)



(iii)



(iv)

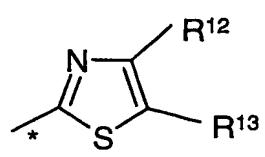
wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R^{10} is at any position on the ring and R^{10} and R^{11} are independently at each instance H, R^a , halogen, -CN, nitro, OR^a , CF_3 , $-NR^aR^a$, $-C(=O)OR^a$, $-C(=O)R^a$, $-C(=O)NR^aR^a$, $-OC(=O)C_{1-4}alkyl$, $-NR^aC(=O)C_{1-4}alkyl$ or $-S(=O)_nR^c$; and wherein R^{11a} is R^a , $-S(=O)_2NR^aR^a$ or $-S(=O)_nR^c$ and $n=1$ or 2.

17. A compound as recited in Claim 14 wherein:

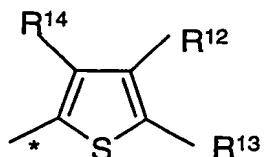
R^4 is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

15

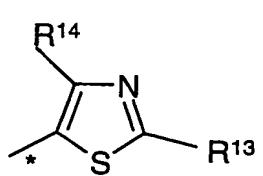
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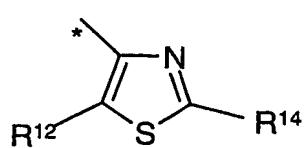
(a)



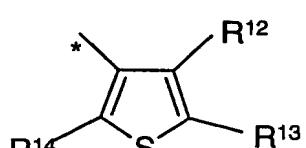
(b)



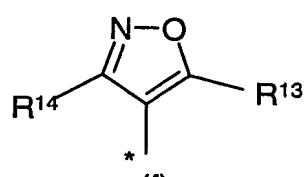
(c)



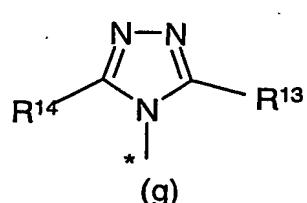
(d)



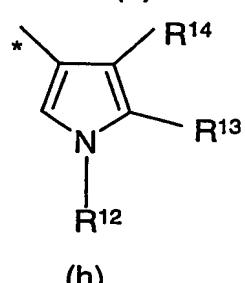
(e)



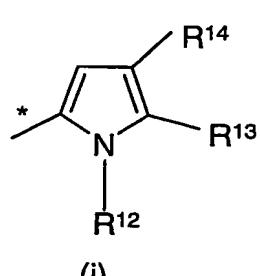
(f)



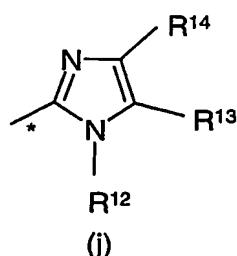
(g)



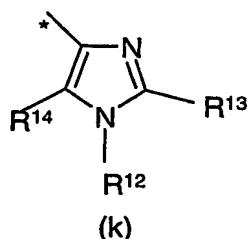
(h)



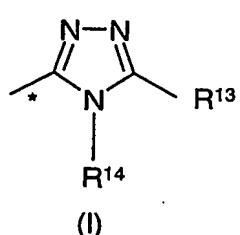
(i)



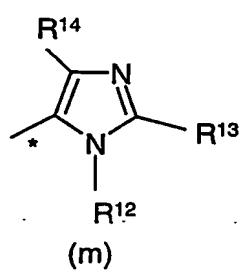
(j)



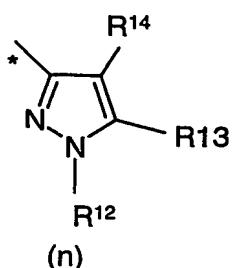
(k)



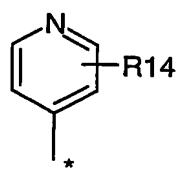
(l)



(m)

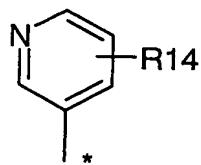


(n)

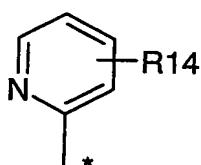


(o)

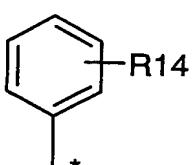
- 108 -



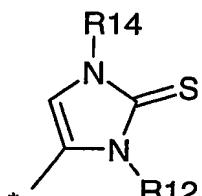
(p)



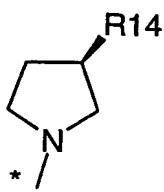
(q)



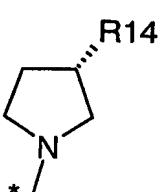
(r)



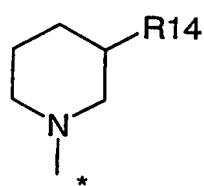
(s)



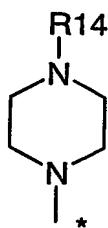
(t)



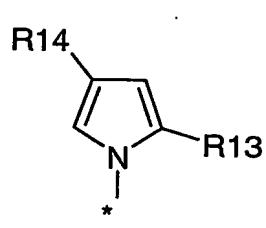
(u)



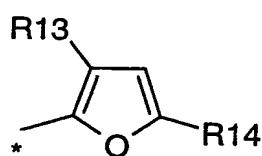
(v)



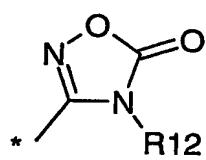
(w)



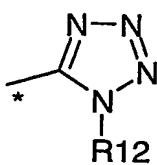
(x)



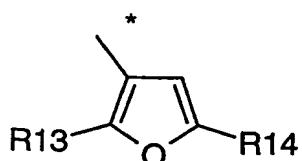
(y)



(z)



(aa)



(ab)

5

wherein * is the location wherein R⁴ is attached to the ring system and wherein
 wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -
 NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -

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C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a).

18. A compound as recited in Claim 14 wherein:

X is S, O, or NR²¹; or XR²⁰ is hydrogen,

W is S, O, or NR²⁰,

10 R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a;

R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a;

R²⁰ and R²¹ and the N to which they are attached in combination can also form a 3 to

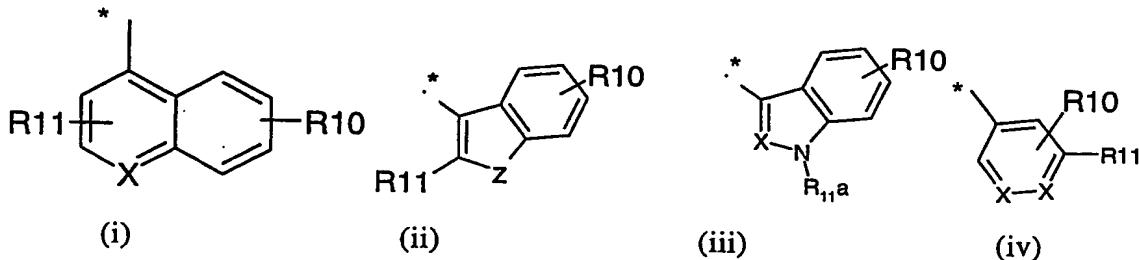
15 10 member N-linked saturated or unsaturated heterocycle having either 1, or 2 heteroatoms independently selected from N, O, or S wherein the heterocycle is substituted with R^e;

10 R¹ is CH₃, CH₂CH₃, CH₂CN, CF₃, (CH₂)₂OH, cyclopropyl, isopropyl, CH₂CCH, (CH₂)₂N(CH₂)₂, (CH₂)₂N(C=NH)NH₂, -CH₂-2-pyridyl, -CH₂-3-pyridyl, -CH₂-4-pyridyl, -(CH₂)₂-1-imidazolyl, -(CH₂)₂-1-pyrazolyl, -(CH₂)₂-1-piperidyl, -(CH₂)_m-(1-methylpiperidin-4-yl), -CH₂-(1-methylpiperidin-3-yl), -(CH₂)₂-(morpholin-4-yl),

20 R² is -CH₂CH₂CH₃, -CH₂-cyclopropyl, -CH₂CH(CH₃)₂, -CH₂CH₂CH₂F, -CH₂-cyclobutyl, -CH₂C(CH₃)₃, -CH₂CH₂CH(CH₃)₂, -CH₂CF₃, -CH₂-methylphenyl, -CH₂-phenol, -CH₂-(3,5-dimethylisoxazol-4-yl), -CH₂-S-phenyl, -CH₂-phenylcarboxyl, or -CH₂SCF₃;

R³ is selected from formulas (i), (ii), (iii) or (iv) set forth below:

25



wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I),

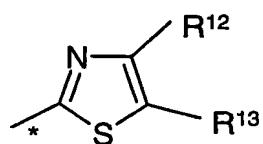
30 and X is C or N; and Z is O or S, wherein R¹⁰ is at any position on the ring and R¹⁰ and R¹¹ are

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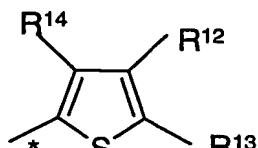
independently at each instance H, R^a, halogen, -CN, nitro, OR^a, CF₃, -NR^aR^a, -C(=O)OR^a, -C(=O)R^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl, -NR^aC(=O)C₁₋₄alkyl or -S(=O)_nR^c; and wherein R^{11a} is R^a, -S(=O)₂NR^aR^a or -S(=O)_nR^c and n=1 or 2.

R⁴ is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

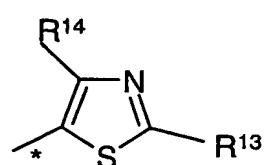
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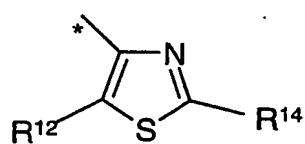
(a)



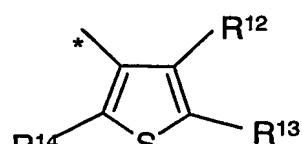
(b)



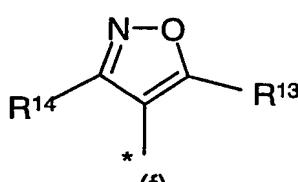
(c)



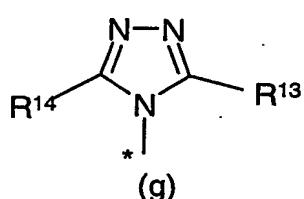
(d)



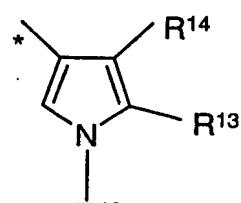
(e)



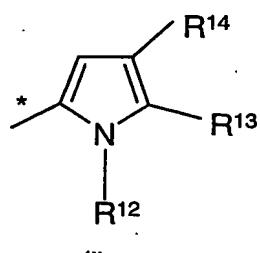
(f)



(g)

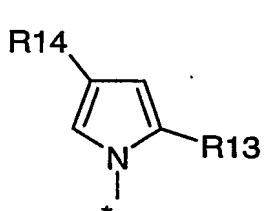
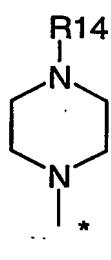
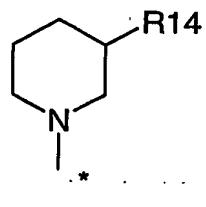
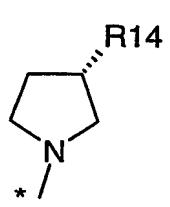
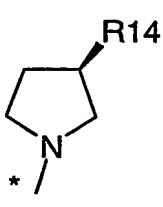
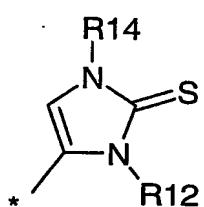
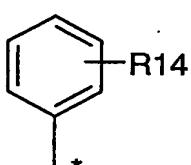
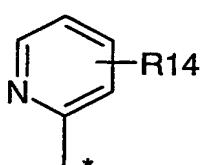
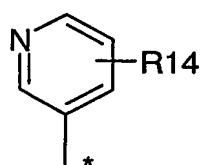
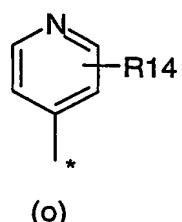
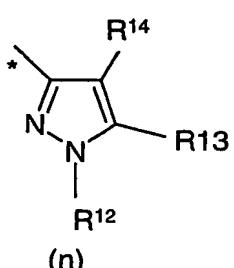
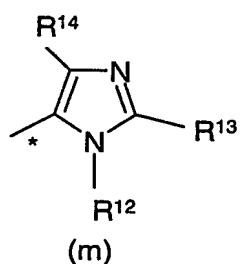
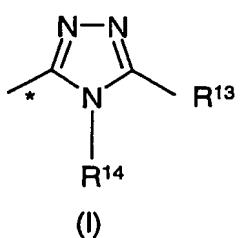
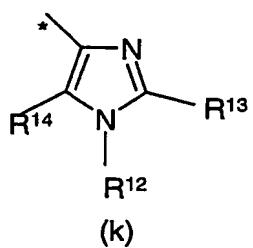
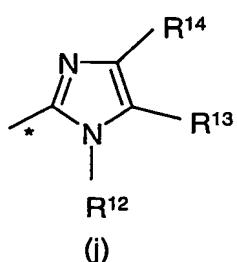


(h)

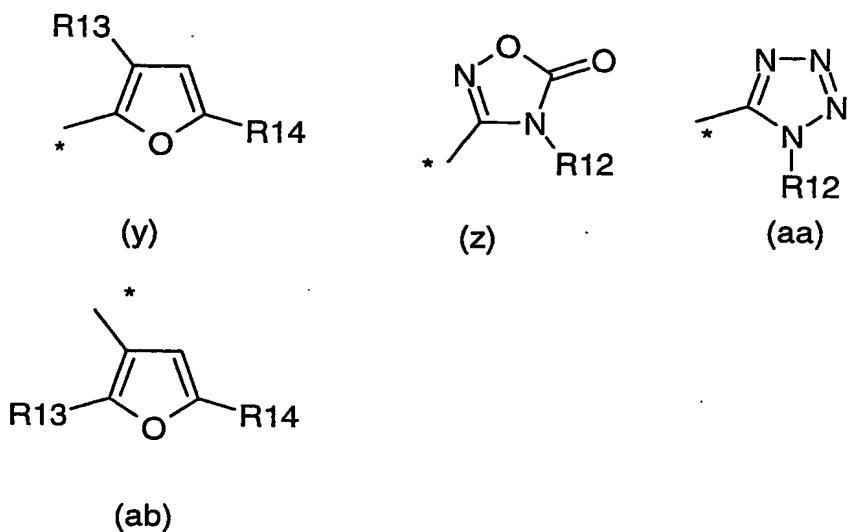


(i)

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- 112 -



wherein * is the location wherein R⁴ is attached to the ring system and wherein
wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -

5 NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -
C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -
C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -
C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -
NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -
10 S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a.

19. A compound of formula (III) selected from:

4-amino-7-isobutyl-2-(1-naphthylmethyl)-3-pyridin-4-yl-2,7-dihydro-6H-pyrazolo[3,4-d]pyrimidin-6-one;

15 7-isobutyl-4-(methylamino)-2-(1-naphthylmethyl)-3-pyridin-4-yl-2,7-dihydro-6H-pyrazolo[3,4-d]pyrimidin-6-one;

4-(dimethylamino)-7-isobutyl-2-(1-naphthylmethyl)-3-pyridin-4-yl-2,7-dihydro-6H-pyrazolo[3,4-d]pyrimidin-6-one;

7-isobutyl-4-(4-methylpiperazin-1-yl)-2-(1-naphthylmethyl)-3-pyridin-4-yl-2,7-dihydro-6H-
20 pyrazolo[3,4-d]pyrimidin-6-one;

4-amino-2-[(6-chloroquinolin-4-yl)methyl]-7-isobutyl-3-(1-methyl-1H-pyrrol-2-yl)-2,7-dihydro-6H-pyrazolo[3,4-d]pyrimidin-6-one;

5-{4-amino-2-[(6-chloroquinolin-4-yl)methyl]-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile;

5-[2-[(6-chloroquinolin-4-yl)methyl]-7-isobutyl-4-(methylamino)-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[2-[(6-chloroquinolin-4-yl)methyl]-4-(dimethylamino)-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[2-[(6-chloroquinolin-4-yl)methyl]-7-isobutyl-6-oxo-4-(propylamino)-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-{2-[(6-chloroquinolin-4-yl)methyl]-4-[(2-hydroxyethyl)amino]-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile

5-[2-[(6-chloroquinolin-4-yl)methyl]-4-(hydroxyamino)-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[2-[(6-chloroquinolin-4-yl)methyl]-4-(cyclopropylamino)-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-{2-[(6-chloroquinolin-4-yl)methyl]-4-hydrazino-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile;

5-[2-[(6-chloroquinolin-4-yl)methyl]-4-(2,2-dimethylhydrazino)-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

N-[2-[(6-chloroquinolin-4-yl)methyl]-3-(4-cyano-1-methyl-1H-pyrrol-2-yl)-7-isobutyl-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-4-yl]acetamide;

5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-4-(methylthio)-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-{2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-4-[(2-hydroxybutyl)amino]-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl}-1-methyl-1H-pyrrole-3-carbonitrile;

5-(2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-4-[(2R)-2-hydroxypropyl]amino)-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-4-methoxy-6-oxo-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

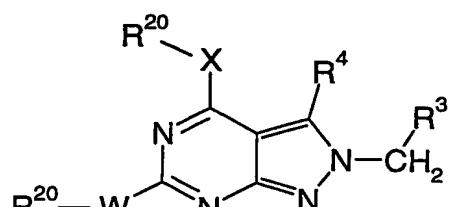
5-[2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-oxo-4-(1H-pyrrol-1-yl)-6,7-dihydro-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1H-pyrrole-3-carbonitrile;

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5-[(6Z)-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-4-(methylamino)-6-(methylimino)-6,7-dihydro-2*H*-pyrazolo[3,4-*d*]pyrimidin-3-yl]-1-methyl-1*H*-pyrrole-3-carbonitrile;

5-4-amino-2-[(6-chloroquinolin-4-yl)methyl]-7-(cyclopropylmethyl)-6-oxo-6,7-dihydro-2*H*-pyrazolo[3,4-*d*]pyrimidin-3-yl]-1-methyl-1*H*-pyrrole-3-carbonitrile.

20. A compound having the structural formula (IV):



(IV)

wherein,

X is S, O, NR²¹; or XR²⁰ is hydrogen;

W is S, O, or NR²¹;

R³ is a monocyclic or bicyclic, saturated or unsaturated, ring system comprising 0, 1, 2

15 or 3 heteroatoms independently selected from N, O, or S, the ring being substituted by 0, 1, 2 or 3 substituents selected from =O, halogen, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, nitro, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -NR^aR^a, -NHC(=O)R^a, N=NR^a,

20 aminocarbonyl, phenyl, benzyl; or R³ is represented by -Het, -Het-Het, R⁵,-R⁵-Het, -Het-R⁵, -Het-O-R⁵, -R⁵-R⁵, -R⁵-OR⁵;

R⁴ is a monocyclic or bicyclic, saturated or unsaturated, ring system, or a vicinal-fused derivative thereof, which may contain from 5 to 12, preferably 5 to 10, ring atoms, 0, 1, 2, 3 or 4 of which are heteroatoms independently selected from N, O, or S, the ring system being substituted by 0, 1, 2 or 3 substituents selected from B(OH)₂, vicinal -OCH₂CH₂O-, vicinal -OC₁₋₂haloalkylO-, vicinal -OCH₂O-, vicinal -CH₂OCH₂O-, =O, halogen, -R^bOR^a, -SR^a, -OR^a, C₁₋₆alkyl, C₁₋₆haloalkyl, -CN, -S(=O)_nR^c, -O(CH₂)_mHet, -O(CH₂)_mC(=O)Het, -

25 O(CH₂)_mC(=O)NR^aR^a, -O(CH₂)_mC(=O)OR^a, -O(CH₂)_mNR^aR^a, -O(CH₂)_mOR^a, -S(CH₂)_mHet, -S(CH₂)_mC(=O)Het, -S(CH₂)_mC(=O)NR^aR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mC(=O)OR^a, -S(CH₂)_mNR^aR^a, -S(CH₂)_mOR^a, -S(CH₂)_mHet, -

$S(CH_2)_mOR^a$, $-NR^aR^a$, $-NHC(=O)R^a$, $-NHC(=O)OR^a$, $N=NR^a$, NO_2 , $-C(=O)NR^aR^a$, $-C(=O)NR^aOR^a$, $-C(=O)NR^a(R^bNR^aR^a)$, $-C(=O)NR^a(R^bOR^a)$, $-C(=O)NR^a(R^bS(=O)_nR^a)$, $-C(=O)NR^a(R^bHet)$, $-C(=O)OR^a$, $-OC(=O)R^a$, $-C(=O)OR^bNR^aR^a$, $-C(=O)R^a$, $-C(=O)R^bNR^aR^a$, $-C(=NOR^a)R^a$, $-C(=NCN)R^a$, $-S(=O)_2NR^aR^a$, $-NR^aS(=O)_2R^a$, $-S(=O)_2NR^a(R^bC(=O)NR^aR^a)$, $-$

5 $S(=O)_2NR^a(R^bC(=O)OR^a)$, aminocarbonyl, phenyl, benzyl; or R^4 is represented by $-(CH_2)_nR^5$ -
Het, $-(CH_2)_nR^d$, -Het, -Het-Het, R^5 , $-R^5$ -Het, -Het- R^5 , -Het- OR^5 , R^5 - R^5 , or $-R^5$ - OR^5 ; or R^4 is
represented by C_{1-6} alkyl, $-NC_{1-6}$ alkyl, or $-N(C_{1-6}$ alkyl)₂ wherein the C_{1-6} alkyl, $-NC_{1-6}$ alkyl,
 $-N(C_{1-6}$ alkyl) are substituted by 0, 1 or 2 substituents selected from R^a , OR^a , halogen or
phenyl wherein R^4 is not $-(CH_2)_2CH_3$, $-(CH_2)_2CH_2OH$, $-(CH_2)_2CO_2H$, or $-(CH_2)_2CO_2C_{1-6}$ alkyl
wherein z is 1,2,3,4,5, or 6;

R^5 is independently at each instance, phenyl substituted by 0, 1, 2, or 3 groups selected
from halogen, C_{1-6} haloalkyl, $-OC_{1-6}$ haloalkyl, C_{1-6} alkyl, $-CN$, nitro, $-OR^a$, $-S(=O)_nR^c$,
 $-O(CH_2)_mHet$, $-O(CH_2)_mC(=O)Het$, $-O(CH_2)_mC(=O)NR^aR^a$, $-O(CH_2)_mC(=O)OR^a$,
 $-O(CH_2)_mNR^aR^a$, $-O(CH_2)_mOR^a$, $-S(CH_2)_mHet$, $-S(CH_2)_mC(=O)Het$, $-S(CH_2)_mC(=O)NR^aR^a$,
15 $-S(CH_2)_mC(=O)OR^a$, $-S(CH_2)_mNR^aR^a$, $-S(CH_2)_mOR^a$, $-R^bOR^a$, $-SR^a$, $-C(=O)NR^aR^a$, $-$
 $C(=O)NR^aOR^a$, $-C(=O)NR^aR^bNR^aR^a$, $-C(=O)NR^aR^bOR^a$, $-C(=O)NR^aR^bS(=O)_nR^a$, $-$
 $C(=O)NR^aR^bHet$, $-C(=O)OR^a$, $-OC(=O)R^a$, $-C(=O)OR^bNR^aR^a$, $-C(=O)R^a$, $-C(=O)R^bNR^aR^a$, $-$
 $C(=NOR^a)R^a$, $-C(=NCN)R^a$, $-S(=O)_2NR^aR^a$, $-NR^aS(=O)_2R^a$, $-S(=O)_2NR^aR^bC(=O)NR^aR^a$, or $-$
 $S(=O)_2NR^aR^bC(=O)OR^a$;

20 R^{20} is, independently at each instance, H, $-CN$, $-S(=O)_nR^c$, $-C(=O)R^a$, $-C(=O)NR^aR^a$,
 $-C(=O)OR^a$, $-NR^aC(=O)R^a$, or $-OC(=O)R^a$, optionally substituted alkyl, optionally substituted
alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted
substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl,
optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle,
25 wherein such substitution is selected from cyclopropyl, halogen, nitro, cyano, hydroxy,
trifluoromethyl, amino, carboxy, carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C_{1-4} alkyl,
 C_{2-4} alkenyl, C_{2-4} alkynyl, C_{1-4} alkoxy, C_{1-4} alkanoyl, C_{1-4} alkanoyloxy, $NH(C_{1-4})$ alkyl),
 $N(C_{1-4})alkyl)_2$, C_{1-4} alkanoylamino, $(C_{1-4}alcanoyl)_2$ amino, $N-(C_{1-4}alkyl)carbamoyl$,
 $N,N-(C_{1-4}alkyl)_2carbamoyl$, $(C_{1-4})S$, $(C_{1-4}alkyl)S(O)$, $(C_{1-4}alkyl)S(O)_2$, $(C_{1-4})alkoxycarbonyl$,
30 $N-(C_{1-4}alkyl)sulfamoyl$, $N,N-C_{1-4}alkyl)sulfamoyl$, $C_{1-4}alkylsulfonylamino$, and heterocyclic;

R^{21} is, independently at each instance, H, $-CN$, $-S(=O)_nR^c$, $-C(=O)R^a$, $-C(=O)NR^aR^a$,
 $-C(=O)OR^a$, $-NR^aC(=O)R^a$, or $-OC(=O)R^a$; optionally substituted alkyl, optionally substituted

alkenyl, optionally substituted alkynyl, optionally substituted cycloalkyl, optionally substituted cycloalkenyl, optionally substituted cycloalkynyl, optionally substituted aryl, optionally substituted alkoxy, optionally substituted amino, optionally substituted heterocycle wherein such substitution is selected from cyclopropyl, halogen, nitro, cyano, hydroxy,

5 trifluoromethyl, amino, carboxy, carboxamido, amidino, carbamoyl, mercapto, sulfamoyl, C₁₋₄ alkyl, C₂₋₄ alkenyl, C₂₋₄ alkynyl, C₁₋₄ alkoxy, C₁₋₄ alkanoyl, C₁₋₄ alkanoyloxy, NH(C₁₋₄ alkyl), N(C₁₋₄ alkyl)₂, C₁₋₄ alkanoylamino, (C₁₋₄ alkanoyl)₂amino, N-(C₁₋₄ alkyl)carbamoyl, N,N-(C₁₋₄ alkyl)₂carbamoyl, (C₁₋₄)S, (C₁₋₄ alkyl)S(O), (C₁₋₄ alkyl)S(O)₂, (C₁₋₄) alkoxycarbonyl, N-(C₁₋₄ alkyl)sulfamoyl, N,N-C₁₋₄ alkyl)sulfamoyl, C₁₋₄ alkylsulfonylamino, and heterocyclic;

10

R²⁰ and R²¹ and the N to which they are attached in combination can also form a 3 to 10 member N-linked saturated or unsaturated heterocycle having either 1, or 2 heteroatoms independently selected from N, O, or S wherein the heterocycle is substituted with R^e;

15 R^a is, independently at each instance, H, C₁₋₆alkyl, -C(=O)C₁₋₄alkyl, C₁₋₄haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

R^b is, independently at each instance, C₁₋₆alkyl, -C(=O)C₁₋₄alkyl, C₁₋₄haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

20 R^c is C₁₋₆alkyl, C₁₋₄haloalkyl, phenyl or benzyl;

R^d is phenyl substituted by 0, 1 or 2 groups selected from -CN, halogen, nitro, C₁₋₆alkyl, C₁₋₄haloalkyl, -OH, -OR^c, -NR^aR^a, -S(=O)_nR^c, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, -OC(=O)R^a, B(OH)₂, vicinyl -OCH₂CH₂O-, vicinyl -OC₁₋₂haloalkylO-, vicinyl -OCH₂O-, vicinyl -CH₂OCH₂O-, phenyl, benzyl and a 5- or 6-membered ring, saturated or unsaturated heterocycle containing 1, 2, 3 or 4 heteroatoms independently selected from N, O, or S;

25 R^e is independently at each instance, H, C₁₋₆alkyl, -C(=O)C₁₋₄alkyl, C₁₋₄haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing 1,2,3, or 4 heteroatoms independently selected from N, O or S;

30 m is 1, 2 or 3;

n is 0, 1 or 2;

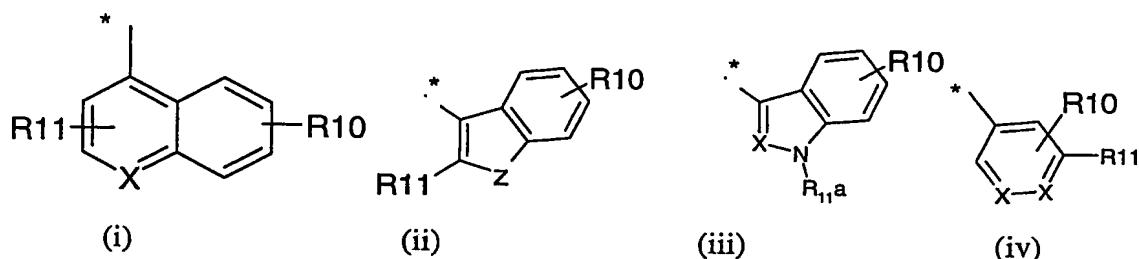
When "optionally substituted" is used, it refers to at least one substituent selected from cyclopropyl, halogen, nitro, cyano, hydroxy, trifluoromethyl, amino, carboxy, carboxamido,

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amidino, carbamoyl, mercapto, sulfamoyl, C₁₋₄ alkyl, C₂₋₄ alkenyl, C₂₋₄ alkynyl, C₁₋₄ alkoxy, C₁₋₄ alkanoyl, C₁₋₄ alkanoyloxy, NH(C₁₋₄ alkyl), N(C₁₋₄ alkyl)₂, C₁₋₄ alkanoylamino, (C₁₋₄ alkanoyl)₂amino, N-(C₁₋₄ alkyl)carbamoyl, N,N-(C₁₋₄ alkyl)₂carbamoyl, (C₁₋₄)S, (C₁₋₄ alkyl)S(O), (C₁₋₄alkyl)S(O)₂, (C₁₋₄) alkoxycarbonyl, N-(C₁₋₄ alkyl)sulfamoyl, N,N-C₁₋₄ alkyl)sulfamoyl, C₁₋₄ alkylsulfonylamino, and heterocyclic
5 or a pharmaceutically acceptable salt thereof.

21. A compound as recited in Claim 20 wherein:

R^3 is selected from formulas (i), (ii), (iii) or (iv) set forth below:

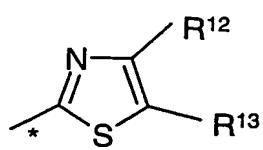


wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R^{10} is at any position on the ring and R^{10} and R^{11} are independently at each instance H, R^a , halogen, -CN, nitro, OR^a , CF_3 , $-NR^aR^a$, $-C(=O)OR^a$, $-C(=O)R^a$, $-C(=O)NR^aR^a$, $-OC(=O)C_{1-4}\text{alkyl}$, $-NR^aC(=O)C_{1-4}\text{alkyl}$ or $-S(=O)_nR^c$; and wherein R^{11a} is R^a , $-S(=O)_2NR^aR^a$ or $-S(=O)_nR^c$ and $n=1$ or 2.

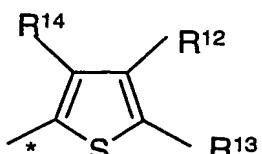
20 22. A compound as recited in Claim 20 wherein:

R^4 is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

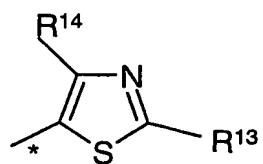
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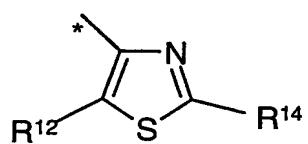
(a)



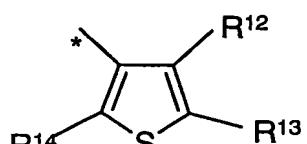
(b)



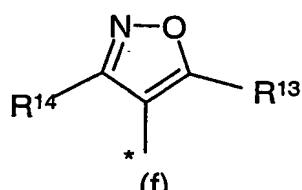
(c)



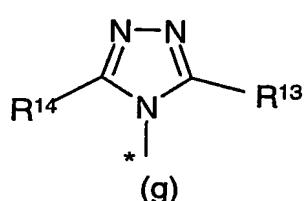
(d)



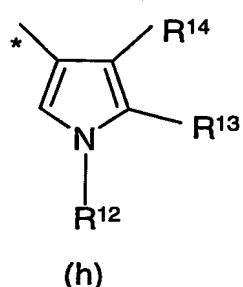
(e)



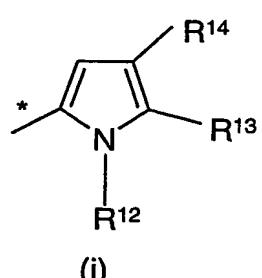
(f)



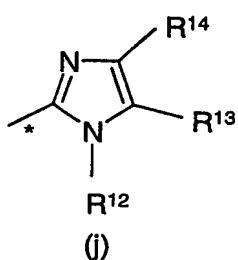
(g)



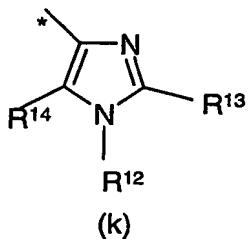
(h)



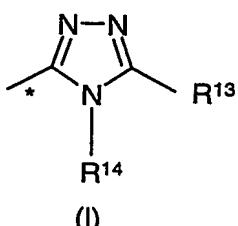
(i)



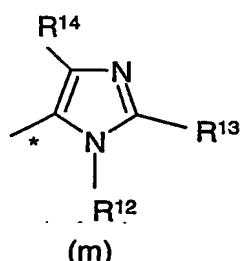
(j)



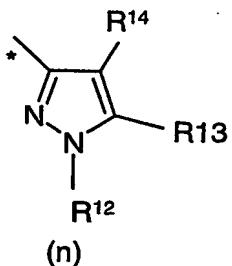
(k)



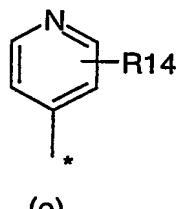
(l)



(m)

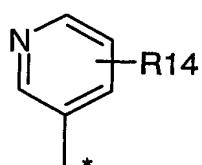


(n)

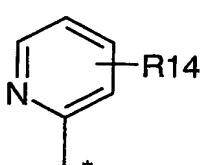


(o)

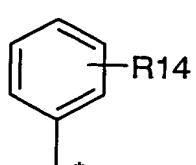
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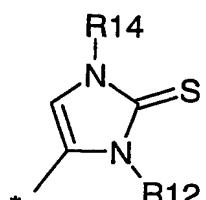
(p)



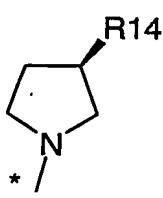
(q)



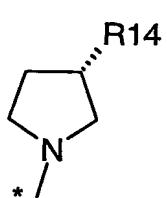
(r)



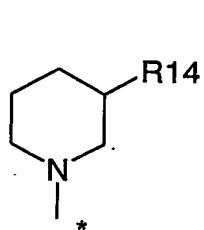
(s)



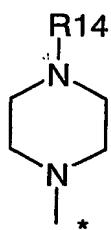
(t)



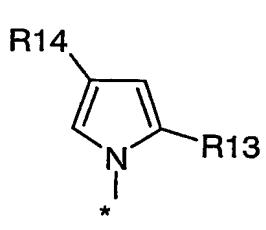
(u)



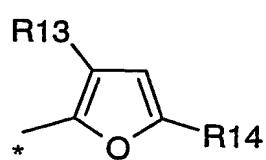
(v)



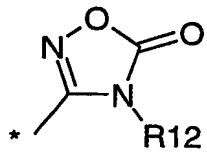
(w)



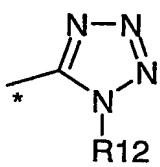
(x)



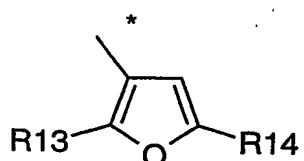
(y)



(z)



(aa)



5

(ab)

wherein * is the location wherein R⁴ is attached to the ring system and wherein
 wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -
 NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)₂R^a, -C(=O)NR^a-Het, -

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C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a).

23. A compound as recited in Claim 20 wherein:

X is S, O, or NR²¹; or X-R²⁰ is hydrogen

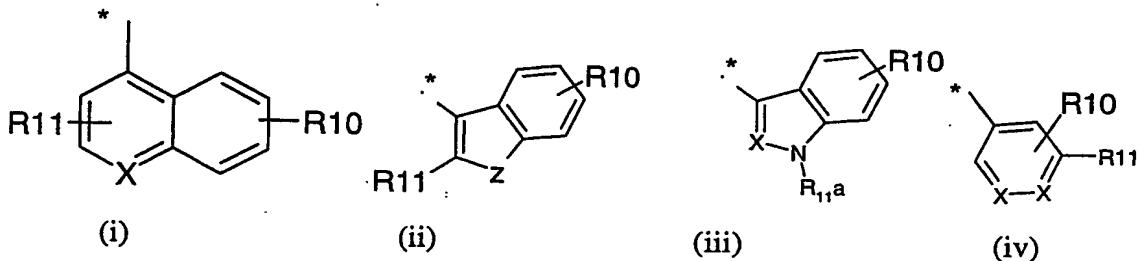
W is S, O, or NR²¹;

10 R¹ is CH₃, CH₂CH₃, CH₂CN, CF₃, (CH₂)₂OH, cyclopropyl, isopropyl, CH₂CCH, (CH₂)₂N(CH₂)₂, (CH₂)₂N(C=NH)NH₂, -CH₂-2-pyridyl, -CH₂-3-pyridyl, -CH₂-4-pyridyl, -(CH₂)₂-1-imidazolyl, -(CH₂)₂-1-pyrazolyl, -(CH₂)₂-1-piperidyl, -(CH₂)_m-(1-methylpiperidin-4-yl), -CH₂-(1-methylpiperidin-3-yl), -(CH₂)₂-(morpholin-4-yl),

R² is -CH₂CH₂CH₃, -CH₂-cyclopropyl, -CH₂CH(CH₃)₂, -CH₂CH₂CH₂F, -CH₂-

15 cyclobutyl, -CH₂C(CH₃)₃, -CH₂CH₂CH(CH₃)₂, -CH₂CF₃, -CH₂-methylphenyl, -CH₂-phenol, -CH₂-(3,5-dimethylisoxazol-4-yl), -CH₂-S-phenyl, -CH₂-phenylcarboxyl, or -CH₂SCF₃;

R³ is selected from formulas (i), (ii), (iii) or (iv) set forth below:

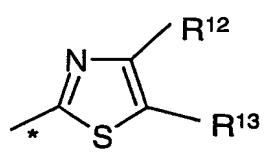


20 (i) (ii) (iii) (iv)

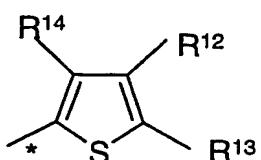
wherein * is the location where (i) or (ii) or (iii) or (iv) is attached to structural formula (I), and X is C or N; and Z is O or S, wherein R¹⁰ is at any position on the ring and R¹⁰ and R¹¹ are independently at each instance H, R^a, halogen, -CN, nitro, OR^a, CF₃, -NR^aR^a, -C(=O)OR^a, -C(=O)R^a, -C(=O)NR^aR^a, -OC(=O)C₁₋₄alkyl, -NR^aC(=O)C₁₋₄alkyl or -S(=O)_nR^c; and wherein R^{11a} is R^a, -S(=O)₂NR^aR^a or -S(=O)_nR^c and n=1 or 2.

R⁴ is selected from formulas (a) to (z) or (aa) or (ab) set forth below:

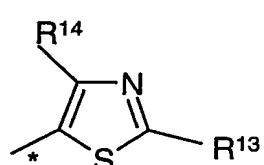
- 121 -



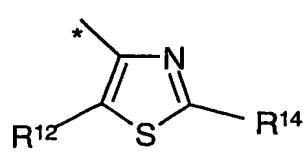
(a)



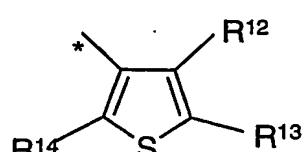
(b)



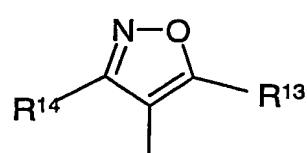
(c)



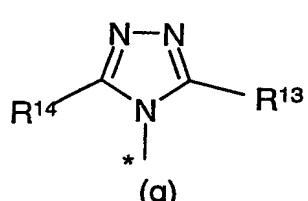
(d)



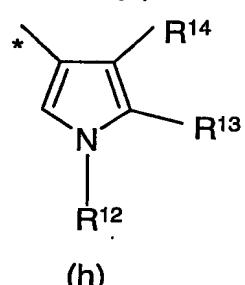
(e)



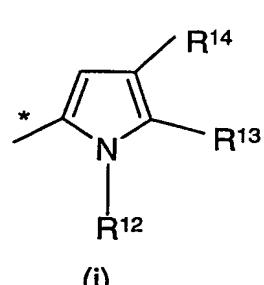
(f)



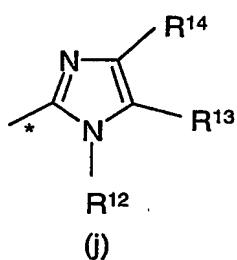
(g)



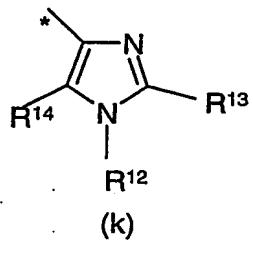
(h)



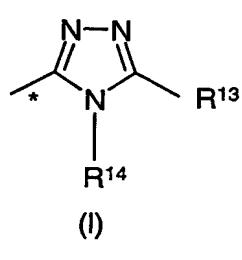
(i)



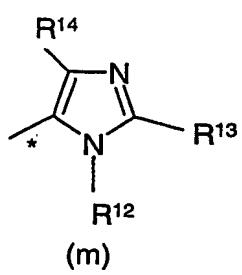
(j)



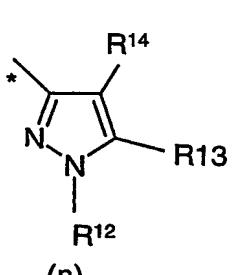
(k)



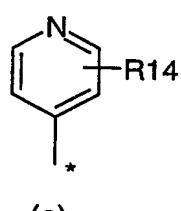
(l)



(m)

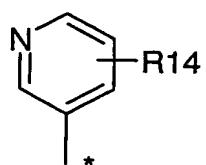


(n)

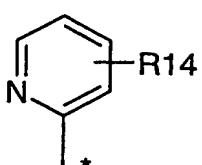


(o)

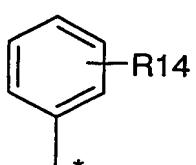
- 122 -



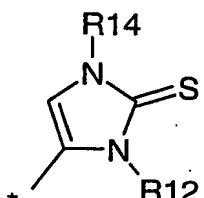
(p)



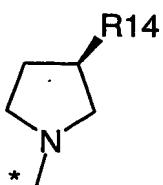
(q)



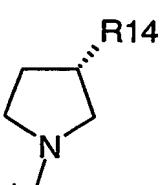
(r)



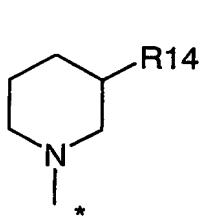
(s)



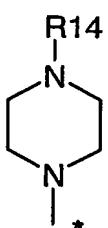
(t)



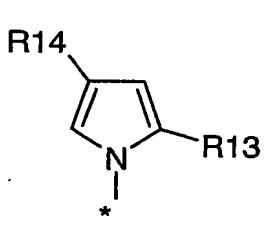
(u)



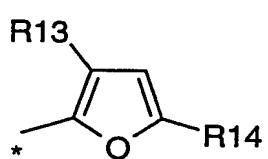
(v)



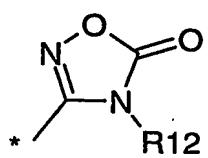
(w)



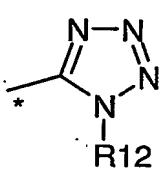
(x)



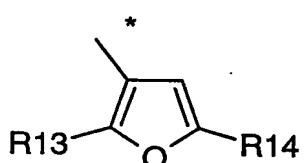
(y)



(z)



(aa)



5

(ab)

wherein * is the location wherein R⁴ is attached to the ring system and wherein
 wherein R¹², R¹³ and R¹⁴ are each independently represented by H, Het, C₁₋₆alkyl, -CN, -NR^aR^a, -nitro, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)NR^aS(=O)R^a, -C(=O)NR^a-Het, -

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C(=O)NR^aNR^aR^a, -C(=O)NR^a(R^bNR^aR^a), -C(=O)NR^a(R^bOR^a), -C(=O)NR^a(R^bS(=O)₂R^a), -C(=O)NR^aR^bHet, -C(=O)NR^aOR^a, -C(=O)R^bNR^aR^a, -C(=NOR^a)R^a, -C(=NCN)R^a, -C(=O)OR^a, -C(=O)OR^bNR^aR^a, -C(=O)R^a, -OC(=O)R^a, -C(=O)R^a-SR^a, =S, -NR^aC(=O)R^a, -NR^aC(=O)OR^a, -NR^aS(=O)₂R^b, -C(=NOR^a)R^a, -S(=O)₂R^a, -S(=O)₂NR^aR^a, -

5 S(=O)₂NR^a(R^bC(=O)NR^aR^a), or -S(=O)₂NR^a(R^bC(=O)OR^a).

R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a;

R²⁰ is H, -CN, R^a, -OR^a, -NR^aR^a, -Het, -S(=O)_nR^c, -C(=O)R^a, -C(=O)NR^aR^a, -C(=O)OR^a, -NR^aC(=O)R^a, or -OC(=O)R^a,

10 R²⁰ and R²¹ and the N to which they are attached in combination can also form a 3 to 10 member N-linked saturated or unsaturated heterocycle having either 1, or 2 heteroatoms independently selected from N, O, or S wherein the heterocycle is substituted with R^e;

R^e is independently at each instance, H, C₁₋₆alkyl, -C(=O)C₁₋₄alkyl, C₁₋₄haloalkyl, phenyl, benzyl, or 5 or 6-membered ring, saturated or unsaturated heterocycle containing

15 1,2,3, or 4 heteroatoms independently selected from N, O or S;

24. A compound of formula (IV) selected from:

5-[2-[(6-chloroquinolin-4-yl)methyl]-6-[(cyclopropylmethyl)amino]-4-(methylamino)-2H-pyrazolo[3,4-d]pyrimidin-3-yl]-1-methyl-1*H*-pyrrole-3-carbonitrile;

20 N-{3-(4-acetyl-1-methyl-1*H*-pyrrol-2-yl)-2-[(6-chloroquinolin-4-yl)methyl]-4-methoxy-2*H*-pyrazolo[3,4-d]pyrimidin-6-yl}-2-cyclopropylacetamide.

25. A compound according to any one of claims 1 to 24, for use as a medicament.

26. The use of a compound as defined in any one of claims 1 to 24, in the manufacture of a medicament for the treatment or prophylaxis of disorders associated with *H. pylori* infection.

27. A method for the treatment of infections associated with *H. pylori* comprising administering to a host in need of such treatment a therapeutically effective amount of a 30 compound as defined in any one of claims 1 to 24.

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28. A method for the prophylaxis treatment of infections associated with *H. pylori* comprising administering to a host in need of such treatment a therapeutically effective amount of a compound as defined in any one of claims 1 to 24.

5 29. A method for the treatment or prophylaxis of *H. pylori* infection comprising administering a therapeutically effective amount of a compound as defined in any one of claims 1 to 24 or a pharmaceutically acceptable salt as claimed in any one of claims 1 to 24.

10 30. A pharmaceutical composition comprising a compound as defined in any one of claims 1 to 24, together with at least one pharmaceutically acceptable carrier, diluent or excipient.